FINAL

BASELINE HUMAN HEALTH RISK ASSESSMENT FOR VOLATILE ORGANIC CHEMICALS IN GROUNDWATER

EKCO HOUSEWARES FACILITY MASSILLON, OHIO

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EXECUTIVE SUMMARY

INTRODUCTION AND OBJECTIVES

A baseline human health risk assessment (HHRA) was conducted for the EKCO Housewares Facility in Massillon, Ohio at the request of EPA Region V (EPA, 1994a). The objective was to assess health risks to a hypothetical future on-site resident from exposure to volatile organic chemicals (VOCs) in the upper and lower groundwater units. The approaches and methodology used were in accordance with U.S. EPA risk assessment guidelines (U.S. EPA, 1989a) and applicable supplemental documents (EPA, 1989b; 1991a; 1991b; 1992a). The final document incorporates revisions that were requested by EPA Region V on 24 October 1994 (EPA, 1994b). Formal responses to these comments are contained in Appendix D. Some specific groundwater issues relating to comments can be found in Appendix C.

The risks presented in this report are upper bound values based on the unlikely event that the EKCO facility will be developed for residential purposes. Moreover, risks determined for the upper unit represent unrealistically high values in view of the low probability that this groundwater, untreated, would ever be used for drinking water purposes or other household uses. The uncertainties are discussed in detail in Section 6.

DATA EVALUATION

Data evaluated for this HHRA consisted of sample measurements from the groundwater wells located in the upper (shallow, intermediate) and lower (bedrock) units.

Tables ES-1 and ES-2 present the frequency of detection, range of detected concentrations, mean concentrations, 95% upper confidence limits (95% UCL) concentrations, and exposure concentrations for each chemical detected in the upper and lower units. Section 2 of this

report presents the detailed evaluation of sample data. Appendix A contains the raw data that were evaluated.

EXPOSURE ASSESSMENT

On-site wells were evaluated for groundwater use by a hypothetical future resident assuming shutdown of the production wells and unrestricted future use of the property. Both upper and lower units were evaluated for the following exposure pathways:

- Drinking water ingestion
- Inhalation while showering
- Dermal contact while showering
- Garden produce ingestion

Guidance for calculating exposure doses (daily intakes) was obtained from U.S. EPA (1989a; 1989b; 1991a; 1992a). Detailed dose information can be found in the tables in Section 5 and Appendix B of this report.

TOXICITY ASSESSMENT

Toxicity criteria for carcinogenic and noncarcinogenic effects from IRIS (U.S. EPA, 1994c) and HEAST (U.S. EPA, 1994d) were used according to U.S. EPA (1989a) guidance. Details of the toxicity assessment are presented in Section 4 of this report.

RISK CHARACTERIZATION

The risk results are provided in Section 5 of this report. Tables ES-3 and ES-4 summarize the cancer and noncancer risks by groundwater unit, exposure pathway and age-group.

Table ES-1

Data Summary for the Shallow Groundwater Unit EKCO Housewares Massillon, Ohio (All Concentrations in µg/L)

Chemical	Frequency of Detection	Range of Detected Concentrations	Mean Concentration*	Upper 95 Percent Confidence Limit Concentration	Exposure Point Concentration	Maximum Contaminant Level ^b
Acetone	3 / 32	4.4 - 17	3.27	4.25	4.25	NE
Benzene	2 / 32	0.15 - 4	1.1	1.45	1.45	5
2-Butanone	1 / 32	2.6	2.22	2.6	2.6	NE
Carbon disulfide	2 / 32	1 - 1.1	1.12	1.1	1.1	NE
Chloroethane	3 / 32	2 - 6.4	2.51	3.19	3.19	NE
Chloroform	3 / 32	1 - 10	1.4	1.99	1.99	100
1,1-Dichloroethane	27 / 32	1 - 2,200	279	451	451	NE
1,2-Dichloroethane	4 / 32	3.7 <i>- 7</i> 3	5	9.24	9.24	5
1,1-Dichloroethene	15 / 32	1.5 - 1,900	109	220	220	7
1,2-Dichloroethene (total)	18 / 32	1 - 480	61.4	93.7	93.7	70⁴
Ethylbenzene	3 / 32	0.22 - 3	1.08	1.41	1.41	700
4-Methyl-2-pentanone	1 / 32	7	2.27	2.92	2.92	NE
Tetrachloroethene	4 / 32	0.86 - 55	4.29	7.71	7.71	5
Toluene	5 / 32	0.25 - 130	5.77	12.6	12.6	1,000
1,1,1-Trichloroethane	18 / 32	2.5 - 52,000	2,490	5,380	5,380	200
1,1,2-Trichloroethane	3 / 32	7.4 - 140	7.58	15.6	15.6	5
Trichloroethene	23 / 32	1 - 220,000	10,100	22,300	22,300	5
Vinyl chloride	12 / 32	4 - 150	18.2	28.4	28.4	2
Xylenes (total)	2 / 22	1.4 - 5	1.19	1.56	1.56	10,000

^a Calculated using "proxy" concentrations for nondetects equal to 1/2 of the sample reporting limit.

^b U.S. EPA Maximum Contaminant Level (MCL) (U.S. EPA, 1994b).

[°] Value is for cis isomer. (MCL for trans isomer is 100 μ g/L.)

NE - An MCL has not been established for this compound.

TABLE ES-2

DATA SUMMARY FOR THE BEDROCK GROUNDWATER UNIT EKCO HOUSEWARES MASSILLON, OHIO

(All Concentrations in $\mu g/L$)

Chemical	Frequency of Detection	Range of Detected Concentrations	Mean Concentration	Upper 95% Confidence Limit Concentration	Exposure Point Concentration	Maximum Contaminant Level ⁶
1,1-Dichloroethane	5/6	2 - 150	52.4	102	102	NE
1,1-Dichloroethene	4/6	1 - 56	13.4	31.3	31.3	7
1,2-Dichloroethene (total)	5/6	4 - 260	59.0	142	142	70⁴
1,1,1-Trichloroethane	4/6	5 - 1,200	216	613	613	200
Trichloroethene	6/6	5 - 430	98.7	234	234	5
Vinyl chloride	2/6	3.3 - 15	3.38	8.15	8.15	2

^{*} Calculated using "proxy" concentrations for nondetects equal to 1/2 of the sample reporting limit.

^b U.S. EPA Maximum Contaminant Level (MCL) (U.S. EPA, 1994b).

 $^{^{\}circ}$ Value is for cis isomer. (MCL for trans isomer is 100 μ g/L.)

NE - An MCL has not been established for this compound.

Table ES-3

Summary of Potential Carcinogenic Health Risks For the Upper and Lower Units by Exposure Pathway and Age Group Future On-Site Resident

EKCO Housewares, Massillon, Ohio

		Upper Unit			Lower Unit	
Exposure Pathway	Individual Cancer Risk	Adult Cancer Risk	Individual Cancer Risk	Child Cancer Risk	Adult Cancer Risk	Individual Cancer Risk
Ingestion of Groundwater	2E-03	4E-03	6E-03	2E-04	4E-04	6E-04
Dermal Absorption while Bathing	1E-04	2E-04	3E-04	7E-06	7E-06	2E-05
Noningestion Groundwater Uses	2E-02	1E-02	3E-02	2E-03	2E-03	3E-03
Ingestion of Garden Produce Irrigated with Groundwater	3E-04	4E-04	7e-04	1E-05	1E-05	3E-05
TOTAL	2E-02	2E-02	4E-02	2E-03	2E-03	4E-03

Carcinogenic Risks

As seen in Table ES-3, total (adult and child) lifetime cancer risks were 4E-02 for the upper unit and 4E-03 for the lower unit. For the upper unit, approximately 90% of the risk was from potable water uses, mainly from inhalation while showering and a smaller portion from water ingestion. Garden produce ingestion and dermal absorption while showering accounted for a small fraction (<3%) of total risk but was nevertheless significant at 1E-03. For the lower unit, cancer risk distribution showed the same pattern as for the upper unit.

Table ES-5 summarizes the cancer risks (1,1-dichloroethane) for the upper and lower units by chemical. The highest contributor overall for the upper unit was 1,1-dichloroethene (1,1-DCE) with about 70% of the total risk. 1,1-DCE accounted for 53% of total risk through the inhalation pathway. Trichloroethene (TCE) was the largest contributor to ridk by the drinking water ingestion pathway. TCE, 1,1-DCE, and vinyl chloride (VC) accounted for the majority of risk through both the garden produce ingestion and dermal absorption pathways.

For the lower unit, 1,1-DCE was the largest contributor to overall risk, with the majority being associated with inhalation while showering, and about 10% that associated with water ingestion. VC and TCE were the second and third highest contributors, respectively, through the same pathways. Dermal absorption while showering risk and garden produce ingestion were the smallest contributors to total risk (about 4E-05 for all three chemicals evaluated).

Noncarcinogenic Risks

Total noncancer risks for the child and the adult are shown by pathway for the upper and lower units in Table ES-4. Total noncancer HIs in the upper unit for all pathways were 420 (child) and 150 (adult). In the lower unit, respective risks were 16 and 4.2. Ingestion of

Table ES-4

Summary of Potential Noncarcinogenic Health Risks for the Upper and Lower Units by Exposure Pathway and Age Group Future On-Site Resident EKCO Housewares Massillon, Ohio

	Uppe	r Unit	Lower Unit		
Exposure Pathway	Child Hazard Index	Adult Hazard Index	Child Hazard Index	Adult Hazard Index	
Ingestion of Groundwater	2.4E+02	1.0E+02	3.8E+00	1.6E+00	
Dermal Absorption while Bathing	1.1E+01	5.9E+00	1.5E-01	8.3E-02	
Noningestion Groundwater Uses (Inhalation)	1.3E+02	2.8E+01	1.1E+01	2.4E+00	
Ingestion of Garden Produce Irrigated with Groundwater	4.4E+01	1.4E+01	5.5E-01	1.7E-01	
TOTAL	4.2E+02	1.5E+02	1.6E+01	4.2E+00	

Table ES-5

Summary of Key Cancer Risks by Chemical in The Upper and Lower Groundwater Units EKCO Facility, Massillon, Ohio

Exposure Pathway	Chemical	Lifetime Cancer Risk
Upper Unit Drinking Water Ingestion	TCE 1,1-DCE VC	4E-03 2E-03 8E-04
Noningestion Uses (inhalation while showering) Noningestion uses (dermal contact while showering)	TCE 1,1-DCE VC 1,1-DCE TCE	1E-03 1E-02 6E-04 1E-04 2E-04
Garden Produce Ingestion	VC TCE 1,2-DCE VC	2E-05 6E-04 1E-04 2E-05
Lower Unit Drinking Water Ingestion	1,1-DCE VC TCE	3E-04 2E-04 1E-04
Noningestion uses (inhalation while showering)	1,1-DCE VC TCE 1,1-DCE	3E-03 2E-04 1E-04
Noningestion uses (dermal contact while showering)	VC TCE	5E-06 2E-06
Garden Produce Ingestion	1,1-DCE VC TCE	1E-05 6E-06 6E-06

TCE = trichloroethene; 1,1-DCE = 1,1-dichloroethene; VC = Vinyl chloride

groundwater in the upper unit provided the highest risks (240 and 100 for the child and adult, respectively). Inhalation while showering provided the next highest risks. Ingestion of garden produce and dermal absorption produced the smallest HIs. For the lower unit, the highest total hazard indices were 11 and 2.4 for the child and adult, respectively through inhalation. Smaller risks were associated with ingestion of drinking water. Again, garden produce ingestion and dermal absorption while showering contributed the least.

Table ES-6 summarizes the noncancer hazard indices for the upper and lower units by chemical. TCE was the largest contributor in the both upper and lower units. In the upper unit, TCE contributed HQs of 240 and 100 for the child and adult, respectively. Garden produce ingestion accounted for 44 and 14, and inhalation accounted for 130 and 28. Dermal contact showed the least risk.

In the lower unit, 1,1-DCE contributed to noncancer risk the greatest with a HQ of 7.6 in the child through inhalation. TCE had HQs of 2.5 and 1.1 for the child and adult, respectively. Dermal absorption and garden produce ingestion were negligible for any individual chemical.

UNCERTAINTY ANALYSIS

Table ES-7 is a summary of uncertainties. These are discussed in greater detail in Section 6 of the report. The greatest uncertainty is the assumption that the upper groundwater unit could be used for future residents for potable purposes (drinking water, showering). This groundwater unit is not currently used off-site as such, except for the potential watering of gardens. Therefore, the risk associated with the upper unit should be completely attributable to garden produce ingestion.

Risks are overestimated in both units because no degradation or dilution of VOCs was assumed over a lifetime of exposure. Normal biodegradation processes and volatilization

Table ES-6

Summary of Key Cancer Risks by Chemical in The Upper and Lower Groundwater Units EKCO Facility, Massillon, Ohio

		Hazard Quotient	
Exposure Pathway	Chemical	Child	Adult
<u>Upper Unit</u> Drinking Water Ingestion	TCE 1,1-DCE VC	240 1.6	110
Noningestion Uses (inhalation while showering)	TCE 1,1-DCE 1,2-DCA	110 12 5	23 2.5 1.1
Noningestion uses (dermal contact while showering) Garden Produce Ingestion	TCE TCE	10 44	5.9
Lower Unit Drinking Water Ingestion	TCE	2.5	1.1
Noningestion (inhalation)	1,2-DCE 1,1-DCE	7.6 1.7	1.6

Table ES-7

Summary of Uncertainty Analysis EKCO Housewares Facility

	Effects on Risk Estimate			
Uncertainty Element	Potential for Overestimation	Potential for Underestimation	Potential for Over- or Underestimation	
Exposure Assessment				
Standard assumptions regarding body weights, skin surface areas, inhalation rates, and life expectancy	i		Low	
Media intake rates	Moderate			
Exposure frequencies	Moderate		1	
Exposure durations	Moderate			
Use of upper unit for residential drinking water and other household uses	High			
Dilution and degradation of VOCs	High			
Toxicity Assessment				
Use of chronic RfDs for estimating noncancer risk in children	Moderate			
Cancer slope (potency) factors	High			
1,1-DichloroethyleneTrichloroethyleneVinyl chloride	Moderate	Moderate	High	
• Reference doses	Moderate-High			

into outdoor air (gardening scenario) will lower the actual exposure point concentrations. Pump and treatment systems in place now or planned for the near future will also substantially reduce potential exposure concentrations over the 30 year future residential lifetime.

CONCLUSIONS

Cancer risks based on future residential development exceed the 1E-04 benchmark used by EPA (1992b) for both upper and lower units, with the upper (4E-02) being one order of magnitude higher than the lower unit (4E-03).

Based on the unlikely potential for the upper unit being used for drinking water and household noningestion uses, and because dilution, degradation and removal were not assumed to occur, the risks for this unit are significantly overestimated and are realistically only related to those exposures that would occur through garden produce ingestion. Risks for the lower unit are also overestimated because degradation, dilution and removal of VOCs was not taken into account during the risk characterization. Finally, based on the low probability that the site will be developed residentially, risks for both units are likely to be hypothetical.

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SECTION 1 INTRODUCTION

1.1 OBJECTIVES

This report is the final baseline human health risk assessment for the EKCO Housewares facility in Massillon, Ohio. A draft of the risk assessment was prepared in response to a request from EPA Region V in a letter dated 22 June 1994 (EPA, 1994a). The final report was prepared after reviewing comments from EPA V as described in their letter of 24 October 1994 (EPA, 1994b). Responses to these comments can be found in Appendix D of this report. The objective is to determine human health risks from exposure to volatile organic chemicals (VOCs) detected in several on-site groundwater units based on the no-action alternative (i.e., in the absence of remedial or corrective action). The no-action alternative is defined for the EKCO facility as the potential future residential development of the site. Note that some interim remediation has been performed for one of the units (the on-site bedrock aquifer), and therefore risks for this unit reflect post-remediation exposure.

The approaches and methodologies used in conducting the baseline risk assessment for the EKCO facility are in accordance with *Risk Assessment Guidance for Superfund* (EPA, 1989) and Supplemental guidance (EPA, 1991a; 1991b, 1991c).

1.2 SITE LOCATION AND DESCRIPTION

1.2.1 Facility Location

The EKCO facility occupies approximately 13 acres located in the town of Massillon, Stark County, Ohio (Figure 1-1). The area immediately surrounding the facility is largely urban and industrial, although land use to the northwest is more rural with a larger proportion of open space. The EKCO property is approximately triangular in shape and is located approximately 1,500 feet west of the Tuscarawas River. The northern border of the facility is Newman Creek, the western border is the Penn Central Railroad, and the eastern border



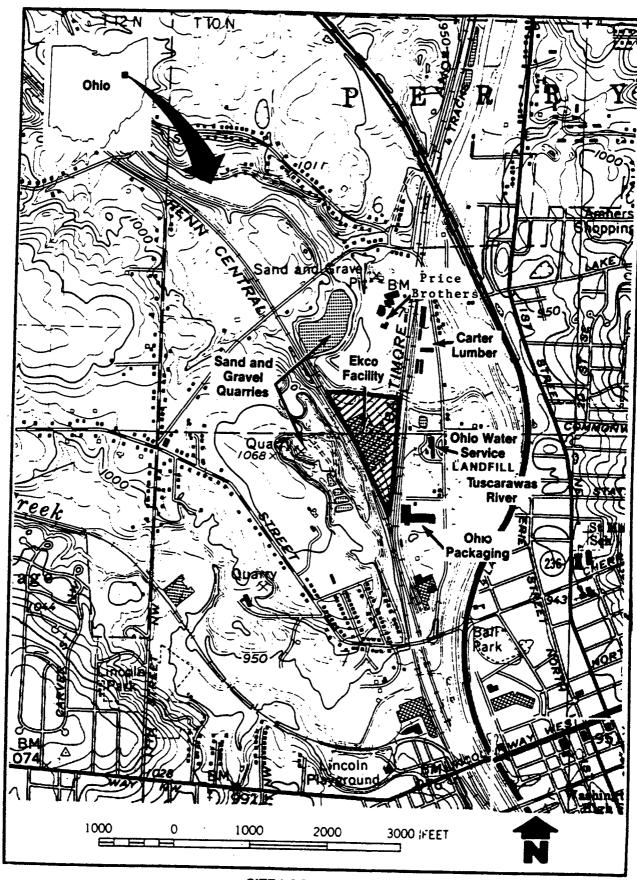


FIGURE 1-1 SITE LOCATION MAP

EKCO HOUSEWARES, INC., MASSILLON, OHIO

(Ref. 7.5 Minute Massillon Quad, Ohio, 1978)

is the Baltimore and Ohio Railroad.

A variety of the business are located adjacent to the EKCO plant. These include Ohio Packaging (paper) to the south, sand and gravel quarries to the west and northwest, Carter Lumber (retail) and Price Brothers (concrete pipe manufacturing) to the north, and the Ohio Water Service (OWS, public water supply) waterworks to the east and northeast. A relatively large, inactive municipal landfill exists just east of the OWS facility. The Baltimore and Ohio Railroad owns numerous spurs and sidings east of and adjacent to the EKCO plant that are used for the storage of rail cars and track maintenance vehicles.

1.2.2 Facility History

A summary of the history of the EKCO facility is presented in Table 1-1. In the 1940s, the EKCO facility manufactured aluminum and stainless steel cookware. By 1951, with the United States becoming involved in the Korean Conflict, the plant began manufacturing 90-mm and 105-mm shell casings for the military. The resulting increase in production necessitated the drilling of two production wells at the facility (W-1 and W-2, as shown in Figure 1-2). In 1953, a sewer was constructed that carried the plant waste to a discharge point along Newman Creek. At approximately the same time, a surface impoundment was constructed along the northern property boundary adjacent to Newman Creek. Sludge from waste treatment was discharged to the surface impoundment.

During 1954, EKCO began coating cookware manufactured at the facility. Chlorinated solvents [primarily trichloroethene (TCE) and 1,1,1-trichloroethane (1,1,1-TCA)] were used to clean the products prior to coating, although these solvents were never used by EKCO at the same time. Between 1954 and 1964, EKCO used TCE for cleaning parts; EKCO changed its process solvent to 1,1,1-TCA in 1964 and has used TCA since then.

In 1965, American Home Products Corporaton (AHPC) acquired EKCO. Porcelain and Teflon coating units were installed at the EKCO facility in 1967. In 1969, with the

TABLE 1-1

EKCO FACILITY HISTORY

Date	EKCO Site History		
Ca 1929-32	First recorded activities at facility. Property is owned by Standard Oil Company.		
Ca 1929-42	Fort Pitt/Massillon Bridge Works - Manufacture of iron and steel bridges and structural iron.		
1945	Manufacturing Aluminum and stainless steel cookware.		
1950	Production wells W-1 and W-2 were installed and put into service to produce water for plant activities. Well W-1 has been used continuously since then, and well W-2 was used until it was taken out of service in the late 1970s.		
1951	The plant began with the U.S. involvement in the Korean conflict manufacturing 90- mm and 105-mm shell casings for the military. This increase in production necessitates the drilling of two production wells (W-1 and W-2).		
1953	A surface impoundment was constructed along the northern property boundary adjacent to Newman Creek, Sludge frame waste treatment was discharged to it. Began copper-plating cookware, used primarily TCE or 1,1,1-TCA to clean cookware.		
1964	Stopped using TCE; 1,1,1-TCA was used in its place.		
1965	AHPC acquired EKCO Housewares.		
1967	Installation of porcelain and teflon coating units.		
1969	Surface impoundment meets newly formed NPDES regulations and permits.		
March 1986	The air stripper system was installed and put into service.		
July 1974	NPDES Permit No. C-3094BD was issued to EKCO.		
1977	EKCO discontinued the manufacturing of aluminum and porcelain cookware and the use of the lagoon ceased.		
1978	All copper plating operations ended; the principal manufactured products were pressed and coated nonstick bakeware.		
1979-1980	The only major documented solvent spill to date at the facility was recorded; neither the exact location nor the extent of the spill was documented.		
1980	The surface impoundments was reactivated under the existing NPDES permit and received housing alkaline degreaser filter water.		
March 1984	In applying for a renewal of their NPDES permit, the plant was required to analyze on-site well water for VOCs, this analysis indicated the presence of 1,1,1-TCA and TCE.		
June 1984	All discharges to lagoon ceased.		
1984	AHPC sold EKCO Housewares to the EKCO Group.		
May 1992	EKCO reported a 330-gallon 1,1,1-TCA spill to EPA. EKCO removed 50 tons of soil from the area of the solvent release.		
Present	EKCO continues to manufacture pressed and coated nonstick bakeware. A silicon-based compound is presently used to coat the bakeware to create the nonstick surface.		

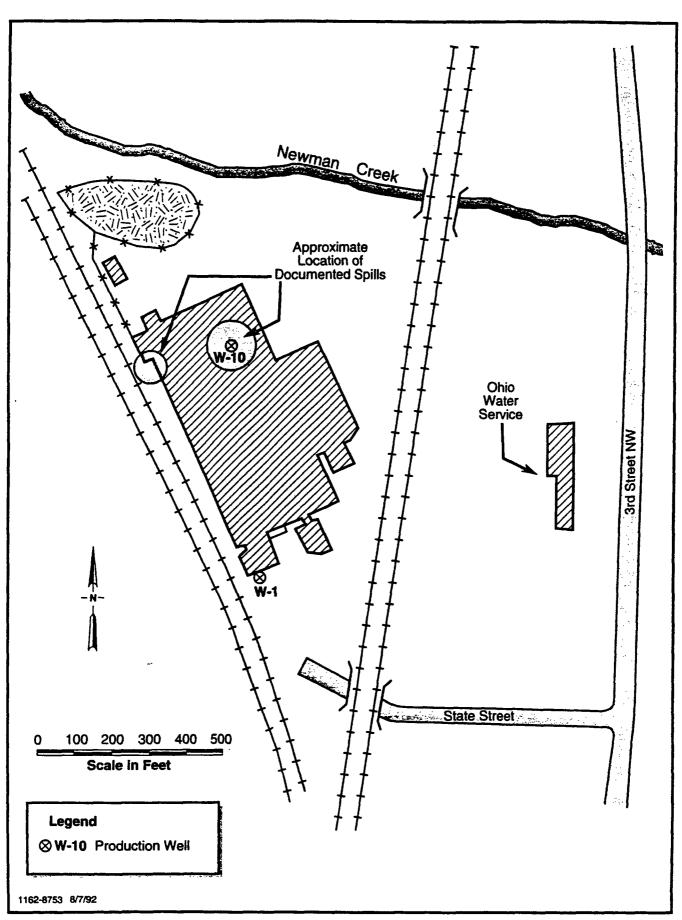


FIGURE 1-2 APPROXIMATE LOCATION OF THE ONLY DOCUMENTED SPILLS

development of the National Pollutant Discharge Elimination System (NPDES) and its attendant regulations and permit requirements, a permit was issued by the State of Ohio for the surface impoundment to discharge waste products associated with plant activities. The waste products permitted for discharge included:

- Deionizers from copper coating operations (hydrochloric acid and sodium hydroxide).
- Washings and waste material from manufacturing porcelain-Teflon coated aluminum cookware (aluminum frit, various pigments: inorganic oxides of lead, cadmium, selenium, and cobalt).
- Alkaline washer fluids to clean aluminum cookware.

In July 1994, NPDES Permit No. C-3094BD was issued to the EKCO facility. As the 1970s progressed, EKCO discontinued the manufacturing of aluminum and porcelain cookware. Use of the surface impoundment ceased in 1977. By the end of 1978, all copper coating operations had ended and the principal products manufactured at the facility became pressed and coated nonstick bakeware.

There have been only two documented solvent releases at the facility. Correspondence between EKCO and the Ohio Environmental Protection Agency (OEPA) identified a solvent release that occurred between 1979 and 1980 in the vicinity of production well W-10 (Figure 1-2). Neither the exact location nor the extent of the spill was documented. Well W-10 is located in a sump covered by a grate mounted flush with the plant floor, and the well head sump may receive floor drainage. In May 1992, EKCO reported to EPA a release of 330 gallons of 1,1,1-TCA in an area northwest of the plant. In response, 50 tons of soil were excavated in the presence of an OEPA representative. The soil was containerized and transported to the Enviro Safe Services of Ohio, Inc. hazardous waste landfill in Toledo, Ohio.

The surface impoundment was reactivated in 1980 under the existing NPDES permit and received alkaline degreaser filter water until mid-1984. In 1984 the facility was acquired by

the EKCO Group, Inc. In March 1984, when the plant applied for a renewal of its NPDES permit, analysis of on-site well water for VOCs was required. The analysis indicated the presence of 1,1,1-TCA and TCE. This discovery resulted in subsequent investigations at EKCO. These investigative activities are described in Subsection 1.3 of this report. The EKCO facility continues to manufacture pressed and coated nonstick bakeware. A silicone-based compound is presently used to coat the bakeware to create the nonstick surface.

1.3 <u>SITE HYDROGEOLOGY</u>

The EKCO facility is situated on the western flank of a glacial valley that extends to the north and south and was carved from Pennsylvanian age sedimentary rocks during Pleistocene glaciation. The majority of the EKCO facility is generally flat except for the northern border of the facility which slopes steeply toward Newman Creek. Surface water runoff at the facility discharges to Newman Creek by two pathways: in the northern part of the facility it flows directly into Newman Creek, and surface water discharge for the rest of the facility is routed through the storm sewer system, which discharges into Newman Creek through Outfall No. 001 located just east of the Baltimore and Ohio Railroad tracks. Figure 1-3 shows the monitor well locations and some important site features.

Prior to the construction of the facility in 1945, a cover of fill material was used to level the natural glacially-formed topography at the building site. Beneath the fill, the glacially deposited sediments form a thin veneer 15 to 30 feet thick in the western portion of the site where the depth to bedrock is shallow. The sediments infill the glacial valley to the east, reaching a maximum on-site thickness of approximately 110 feet at the eastern property boundary. Further off-site to the east, these unconsolidated sediments reach thicknesses exceeding 252 feet.

Figures 1-4 and 1-5 show generalized geologic cross-sections for the stratigraphic units in the area of the site. These cross-sections show that two unconsolidated sand units exist on-site -- the shallow and intermediate units. Off-site, a third deep sand and gravel unit



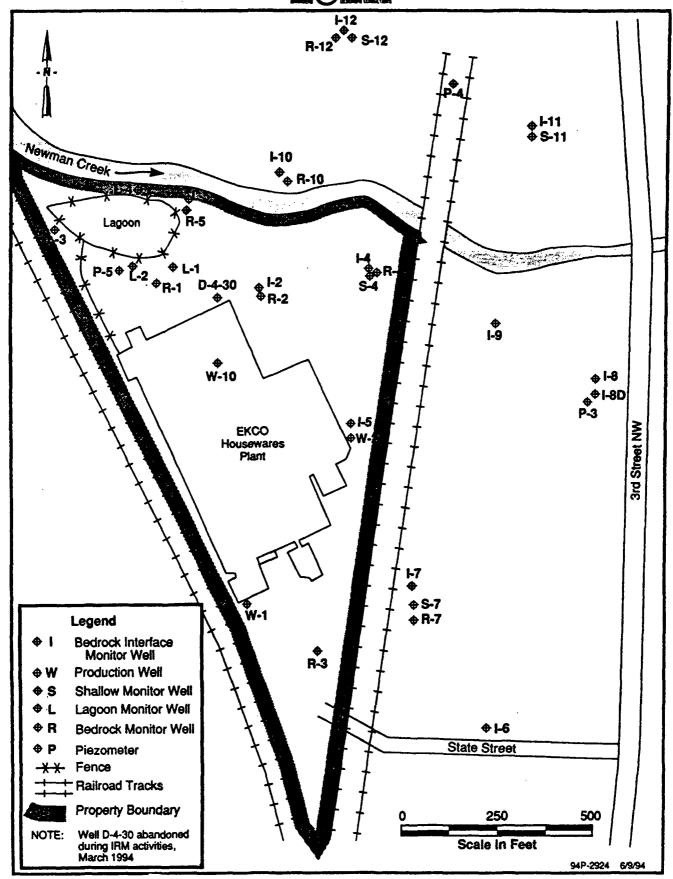
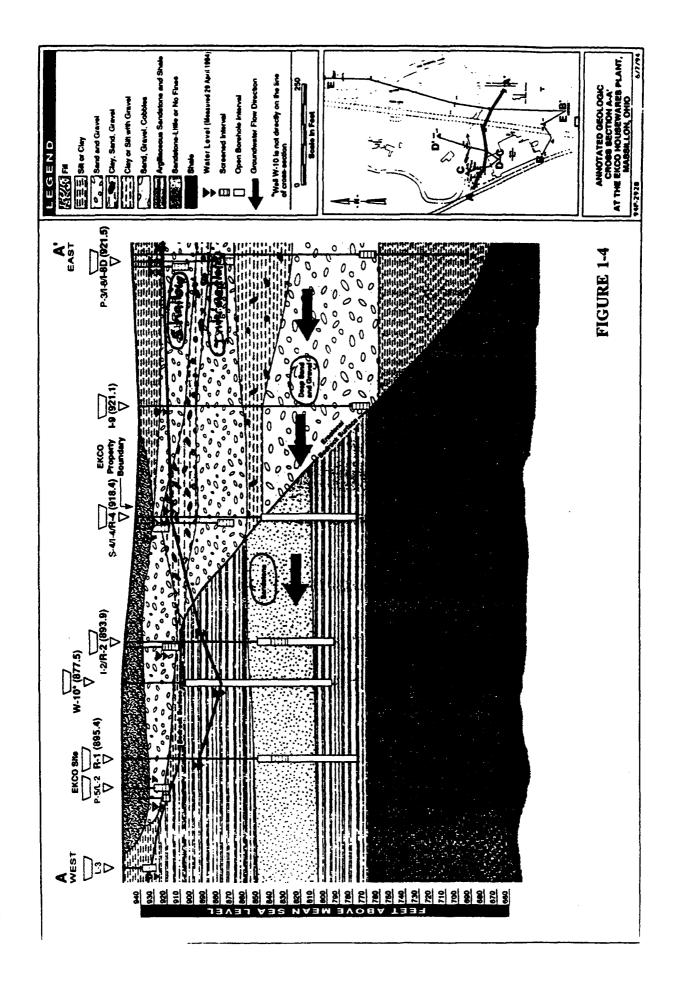
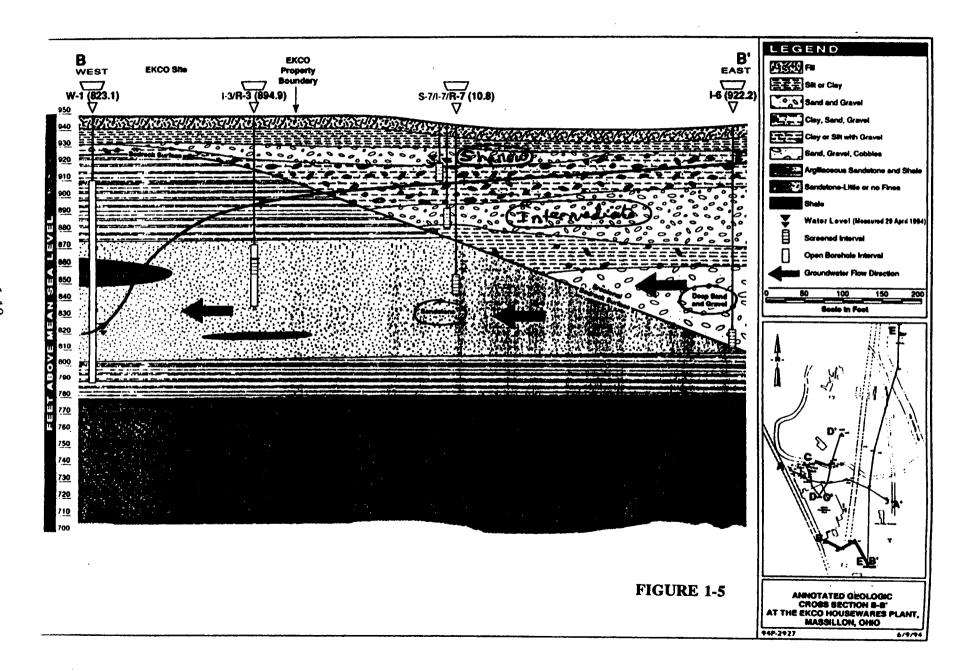


FIGURE 1-3 SITE PLAN EKCO HOUSEWARES FACILITY MASSILLON, OHIO





exists further to the east. These unconsolidated sand units are typically separated by silt and clay units. In general, the shallow, intermediate and dep sand units act as the primary medium for groundwater flow and the low permeable silt and clay units act as barriers to groundwater flow; however, variations in the permeability occur locally. North of well I-2 the silt and clay layer between the shallow and intermediate unit pinches out. In this area the shallow and intermediate units are hydraulically connected to each other.

The bedrock at the site consists of a relatively thick sandstone unit which is both overlain and underlain by layers of shale and argillaceous sandstone. The sandstone unit has a relatively high permeability and is the primary bedrock water bearing unit at the site. The layers of shale and argillaceous sandstone have a relatively low permeability and generally act as barriers to groundwater flow.

There are five groundwater production wells in the area of the site, which have an effect on the local groundwater flow system. EKCO uses two sandstone bedrock production wells W-1 and W-10 pumping at a total of approximately 600 gpm to provide water for the manufacturing facility. Ohio Water Service (OWS) pumps three production wells (OWS-1, -1 and -3) intermittently from the deep sand and gravel unit at a total of up to 2,800 gpm to provide water for the City of Massillon. The OWS production wells are located approximately 2,000 feet northeast of the EKCO facility in the deep sand and gravel glacial valley.

Groundwater contour maps for the shallow and intermediate units are presented on Figures 1-6 and 1-7. The figures show that the groundwater in these two units flows toward the production wells W-1 and W-10. Figure 1-5 shows that the shallow unit is dewatered beneath a large portion of the site, and the cross-sections (Figures 1-4 and 1-5) show that the intermediate unit only exists beneath the eastern portion of the site in the area of wells (I-4 and I-7. The shallow and intermediate units have been grouped together for the risk assessment evaluation because of several similar characteristics. Groundwater yield is not sufficient to support drinking water use from either of these units (See Appendix C). They



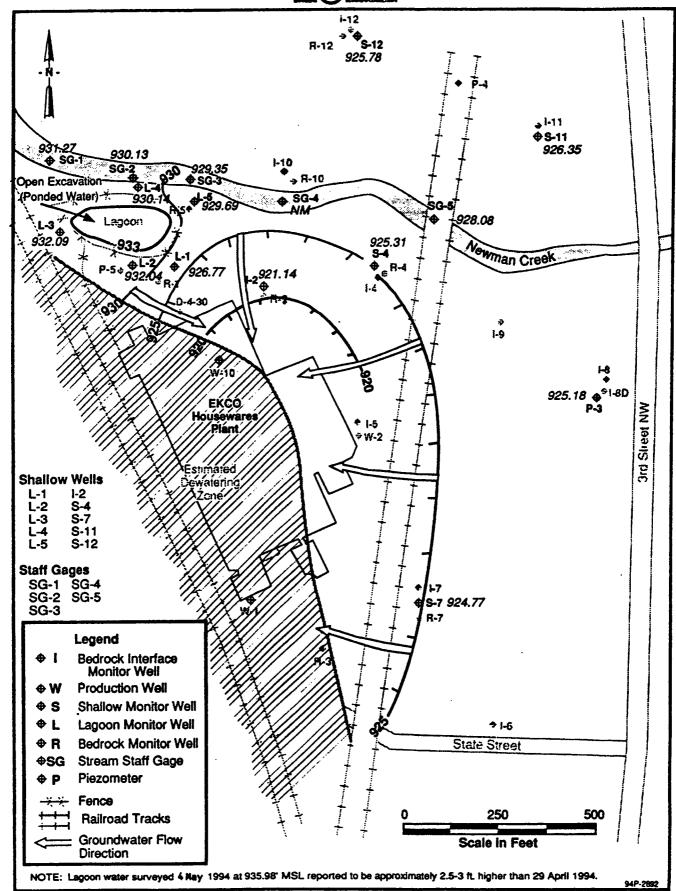
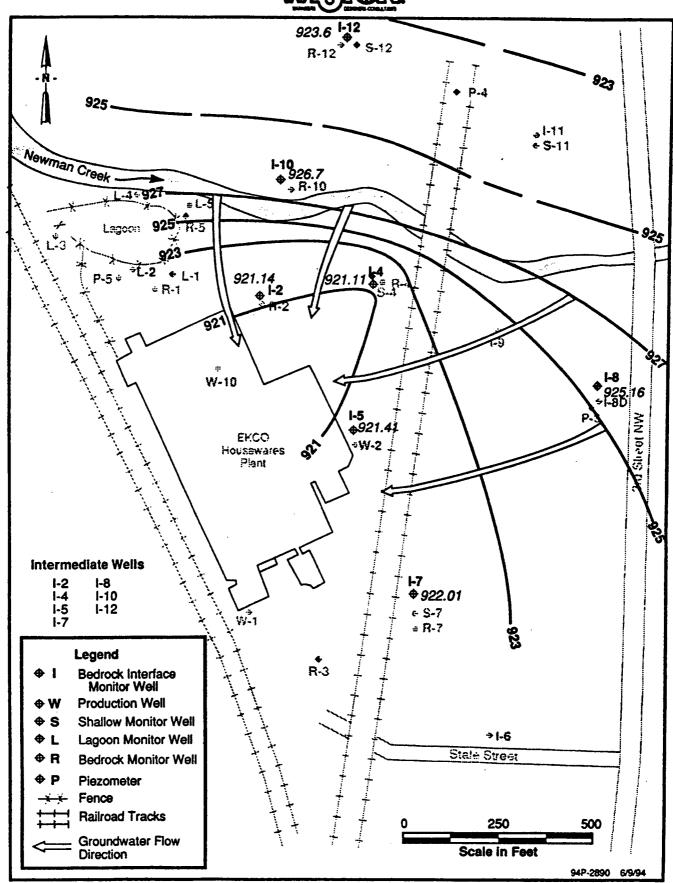


FIGURE 1-6 GROUNDWATER CONTOUR MAP OF WELLS COMPLETED IN THE SHALLOW SAND AND GRAVEL WATER BEARING ZONE - 29 APRIL 1994 EKCO HOUSEWARE FACILITY - MASSILLON, OHIO

WESTERN.



GROUNDWATER CONTOUR MAP OF WELLS COMPLETED IN THE INTERMEDIATE SAND AND GRAVEL WATER BEARING ZONE, WATER LEVELS MEASURED 29 APRIL 1994 EKCO FACILITY – MASSILLON, OHO

only exist in saturated conditions beneath a small portion of site, and geologic logs from the northern portion of the site indicate that these units are hydraulically connected in some areas.

A bedrock groundwater contour map is presented in Figure 1-8 and shows that a significant bedrock drawdown cone exists beneath the entire EKCO site. This drawdown cone is induced by the pumping of EKCO production wells W-1 and W-10. The water from these two production wells is treated by an on-site air stripper system. Off-site residential wells obtain drinking water either from the sandstone bedrock or the deep glacial sand and gravel aquifer.

Packer testing activities conducted during the RCRA Feasibility Investigation (RFI) indicated that some of the bedrock wells may have had leaking casings, allowing groundwater to flow through the well annulus form the shallow overburden to the sandstone bedrock. The casings for bedrock wells R-1, R-2, R-3, W-1, W-2 and W-10 were sealed during Interim Remedial Measure (IRM) activities conducted in April 1994 (WESTON, 1994).

In summary, the risk assessment will evaluate groundwater as two distinct water-bearing zones at the site; the Upper Unit and the Lower Unit. Groundwater within the Upper Unit typically occurs under unconfined or semiconfined conditions and is represented by shallow and intermediate well data. The general direction of groundwater flow in the Upper Unit is toward wells W-1 and W-10. Groundwater within the Lower Unit is typically semiconfined by overlying shale and is represented by on-site bedrock well data. The direction of groundwater flow in the Lower Unit is also toward wells W-1 and W-10.



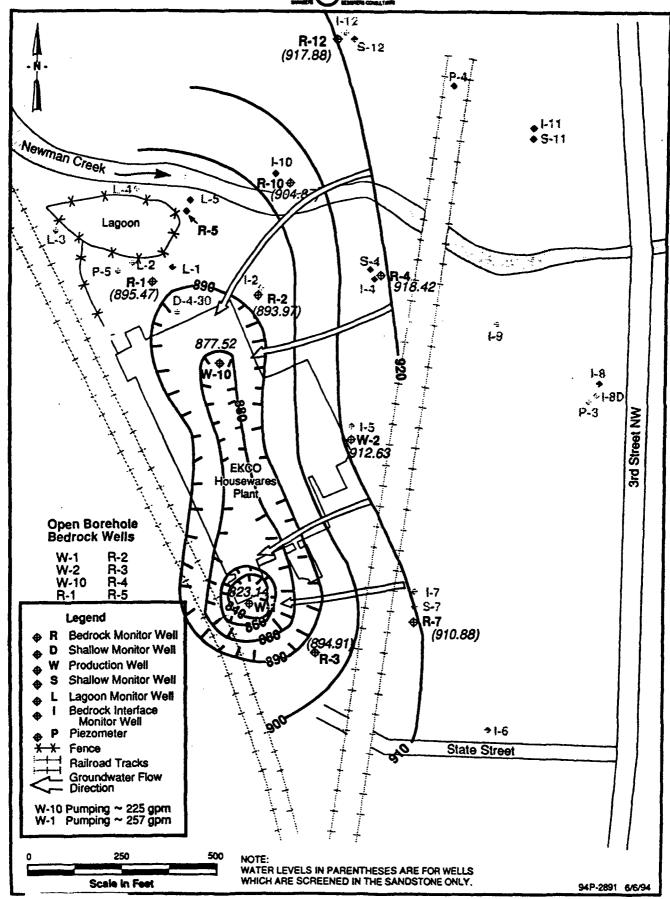


FIGURE 1-8 GROUNDWATER CONTOUR MAP OF WELLS COMPLETED IN THE BEDROCK, 29 APRIL 1994 EKCO HOUSEWARES FACILITY – MASILLON, OHIO

1.4 ORGANIZATION OF REPORT

The remainder of this report is divided into the following sections:

- Data Evaluation (Section 2) presents an evaluation and summary of the chemicals detected in the groundwater units at the EKCO facility. This characterization is comprised of three principal technical elements:
 - The evaluation of groundwater data used in the human health risk assessment.
 - The selection and identification of chemicals of potential concern for which human health risks were assessed.
 - Comparison of the concentrations for the chemicals of potential concern with applicable or relevant and appropriate requirements (ARARs) such as MCLs.
- The Exposure Assessment (Section 3) presents the pathways by which the hypothetical on-site future resident (child and adult) could come in contact with the chemicals of potential concern, the exposure algorithms and input assumptions used, and the calculated exposure concentrations and estimated daily intakes (doses). Guidance for estimating the reasonable maximum exposure (RME) was obtained from EPA (1989).
- The Toxicity Assessment (Section 4) presents a discussion of the carcinogenic and noncarcinogenic toxicity criteria used to evaluate pathway specific human health risk.
- The Risk Characterization (Section 5) summarizes and discusses the risk results. This section will present the risk summary tables.
- The Uncertainty Analysis (Section 6) discusses those chemical/pathway specific risks that had the greatest influence on total risk and the magnitude of overor underestimation of risk that occurred as a result of the assumptions used in the risk assessment. Also included in this section is a discussion of central tendency exposures (EPA, 1992).
- Appendices present (A) the raw data, (B) the exposure dose and risk calculations, (C) groundwater data analysis issues, and (D) WESTON responses to EPA Region V comments on the draft risk assessment.

1.5 REFERENCES

- EPA (U.S. Environmental Protection Agency), 1989. Risk Assessment Guidance for Superfund. Volume I: Human Health Evaluation Manual, Part A. Interim Final. Office of Solid Waste and Emergency Response. Washington, D.C. OSWER Directive 9285.701A.
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- EPA (U.S. Environmental Protection Agency), 1994b. Letter to P. McDonald (American Home Products Corporation, Madison, NJ 17940) from S. Averill (U.S. EPA Region V, RCRA Enforcement Branch, Chicago, IL 60604-3590). 24 October 1994.

SECTION 2 DATA EVALUATION

2.1 APPROACH

The objectives of the data evaluation process were to review and summarize the analytical data for each of the groundwater units examined in the risk assessment in order to select chemicals of potential concern and develop exposure point concentrations. Only on-site well locations were evaluated since only future on-site uses were assessed (see Section 3). Moreover, only volatile organic chemical (VOC) data for the on-site wells were assessed, as requested by EPA (1994a).

2.1.1 Upper Unit

The upper (shallow/intermediate) groundwater unit was characterized using data from sampling events in 1988, 1991, and 1992 for the following wells (analytical data is contained in Appendix A):

- S-4, S-7 (note: no 1988 data available for S-4)
- D-4-30
- L-1, L-2, L-4, L-5 (L-3, background)
- I-2, I-4, I-5, I-7

Wells I-8, I-9, I-11, I-12, I-13, R-12, S-11 and S-12 were at off-site locations and therefore were not evaluated.

2.1.2 Lower Unit

Data from the lower (bedrock) groundwater unit were evaluated using currently available post-Interim Remedial Measures (post-IRM) data (Spring, 1994) for the following on-site wells:

• R-1, R-2, R-3, R-5, W-1, W-10

mand here be wells

Bedrock well data from the years prior to interim remediation were not evaluated since these data no longer represent baseline conditions. Prior to retrofitting, the R-well data represented a combination of the overburden and bedrock groundwater. Retrofitting prevented mixing of the overburden and bedrock groundwater.

2.2 DATA EVALUATION PROCESS

The following guidelines for data evaluation were used to produce the data summaries for each groundwater unit. These approaches are consistent with current U.S. EPA guidelines (U.S. EPA, 1989).

- If a chemical was not positively identified in any sample from a given medium, either because it was not found at a concentration exceeding the sample detection limit or due to blank contamination, it was not addressed for that medium.
- All "J"-qualified data (i.e., estimated concentrations) were used in the quantitative analysis in the same manner as unqualified positive data.
- If a chemical was reported as a non-detect in a particular sample, but was detected in the medium as a whole, one-half of the sample quantitation limit was used as a "proxy" concentration for the chemical in that sample when calculating the mean and 95% upper confidence limit (UCL) concentrations.
- Duplicate sample results from the same sampling round were treated as one data point in summarizing the analytical results. Positive duplicate values were averaged to obtain the single data point; however, if the chemical was detected in one of the duplicates but not the other, the detected concentration was used.
- If a chemical was detected in an investigative sample and was also detected in a field, trip, or laboratory blank associated with that sample, then the sample concentration was considered to be positive only if it exceeded 5 times the blank concentration (or 10 times for the common laboratory contaminants acetone, methylene chloride, toluene, and 2-butanone).
- For the shallow groundwater unit, results for the same well from different sampling rounds were treated as distinct data points.

Tables 2-1 and 2-2 present data summaries for the upper and lower units, respectively. These summaries include the frequency of detection, the range of detected concentrations, the arithmetic mean concentration, and the 95 percent UCL of the mean.

Table 2-1

Data Summary for the Shallow Groundwater Unit EKCO Housewares Massillon, Ohio (All Concentrations in µg/L)

Chemical	Frequency of Detection	Range of Detected Concentrations	Mean Concentration	Upper 95 Percent Confidence Limit Concentration	Exposure Point Concentration	Maximum Contaminant Level*
Acetone	3 / 32	4.4 - 17	3.27	4.25	4.25	NE
Benzene	2 / 32	0.15 - 4	1.1	1.45	1.45	5
2-Butanone	1 / 32	2.6	2.22	2.6	2.6	NE
Carbon disulfide	2 / 32	1 - 1.1	1.12	1.1	1.1	NE
Chloroethane	3 / 32	2 - 6.4	2.51	3.19	3.19	NE
Chloroform	3 / 32	1 - 10	1.4	1.99	1.99	100
1,1-Dichloroethane	27 / 32	1 - 2,200	279	451	451	NE
1,2-Dichloroethane	4 / 32	3.7 - 73	5	9.24	9.24	5
1,1-Dichloroethene	15 / 32	1.5 - 1,900	109	220	220	7
1,2-Dichloroethene (total)	18 / 32	1 - 480	61.4	93.7	93.7	70°
Ethylbenzene	3 / 32	0.22 - 3	1.08	1.41	1.41	700
4-Methyl-2-pentanone	1 / 32	7	2.27	2.92	2.92	NE
Tetrachloroethene	4 / 32	0.86 - 55	4.29	7.71	7.71	5
Toluene	5 / 32	0.25 - 130	5.77	12.6	12.6	1,000
1,1,1-Trichloroethane	18 / 32	2.5 - 52,000	2,490	5,380	5,380	200
1,1,2-Trichloroethane	3 / 32	7.4 - 140	7.58	15.6	15.6	5
Trichloroethene	23 / 32	1 - 220,000	10,100	22,300	22,300	5
Vinyl chloride	12 / 32	4 - 150	18.2	28.4	28.4	2
Xylenes (total)	2 / 22	1.4 - 5	1.19	1.56	1.56	10,000

^a Calculated using "proxy" concentrations for nondetects equal to 1/2 of the sample reporting limit.

^b U.S. EPA Maximum Contaminant Level (MCL) (U.S. EPA, 1994b).

 $^{^{\}circ}$ Value is for cis isomer. (MCL for trans isomer is 100 μ g/L.)

NE - An MCL has not been established for this compound.

TABLE 2-2

DATA SUMMARY FOR THE BEDROCK GROUNDWATER UNIT EKCO HOUSEWARES MASSILLON, OHIO

(All Concentrations in $\mu g/L$)

Chemical	Frequency of Detection	Range of Detected Concentrations	Mean Concentration	Upper 95% Confidence Limit Concentration	Exposure Point Concentration	Maximum Contaminant Level ^t
1,1-Dichloroethane	5/6	2 - 150	52.4	102	102	NE
1,1-Dichloroethene	4/6	1 - 56	13.4	31.3	31.3	7
1,2-Dichloroethene (total)	5 / 6	4 - 260	59.0	142	142	70⁴
1,1,1-Trichloroethane	4/6	5 - 1,200	216	613	613	200
Trichloroethene	6/6	5 - 430	98.7	234	234	5
Vinyl chloride	2/6	3.3 - 15	3.38	8.15	8.15	2

^{*} Calculated using "proxy" concentrations for nondetects equal to 1/2 of the sample reporting limit.

^b U.S. EPA Maximum Contaminant Level (MCL) (U.S. EPA, 1994b).

 $^{^{\}circ}$ Value is for cis isomer. (MCL for trans isomer is 100 μ g/L.)

NE - An MCL has not been established for this compound.

2.3 CHEMICALS OF POTENTIAL CONCERN

The chemicals of potential concern in a risk assessment are those chemicals that are potentially site-related and whose data are of sufficient quality for use in the quantitative analysis. The VOC data for the wells and sampling events listed in Subsection 2.1 were evaluated and summarized according to the guidelines described in Subsection 2.2 above. All VOCs that were positively identified in each groundwater unit after the data evaluation were included as chemicals of potential concern for that unit. Well L-3 served as a background well for the study; however, no VOCs were detected in this well during the 1988, 1991, and 1992 sampling events. Tables 2-1 and 2-2 present the chemicals of potential concern for the upper and lower groundwater units, respectively.

2.4 EXPOSURE POINT CONCENTRATIONS

The exposure point concentration is the concentration of a chemical to which a receptor is expected to be exposed. The exposure point concentration for each chemical of concern in each groundwater unit was calculated as the lower of either the maximum detected concentration or the 95% UCL of the arithmetic mean concentration. The 95% UCL was calculated according to the following standard equation (Mendenhall and Sincich, 1988):

95% UCL =
$$\overline{y} \pm t_{\alpha=0.05} \left[\frac{s}{\sqrt{n}} \right]$$

where:

 $\bar{y} = sample mean$

 $t_{\alpha=0.05}$ = student's *t*-statistic based on (n-1) degrees of freedom

s = sample standard deviation

n = sample size

Exposure point concentrations for the chemicals of potential concern in each groundwater unit are listed in Tables 2-1 and 2-2.

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SECTION 3

EXPOSURE ASSESSMENT

The purpose of the exposure assessment is to estimate the magnitude of human exposure to the chemicals of potential concern at the EKCO facility. The results of the exposure assessment are combined with chemical-specific toxicity information to quantitatively estimate the human health risks associated with chemical exposure (EPA, 1989a).

The identification of actual or potential pathways through which human receptors could be exposed to chemicals in the groundwater at the EKCO facility consists of an evaluation of the contaminated environmental media and a characterization of the potentially exposed populations. After the exposure pathways have been identified, daily intakes (doses) of the chemicals of potential concern can be quantified using standard exposure algorithms (EPA, 1989a).

3.1 PATHWAYS OF EXPOSURE

An exposure pathway generally consists of four elements: (1) a source and mechanism of contaminant release, (2) a retention or transport medium, (3) a point of potential human contact with the contaminated medium (referred to as the exposure point), and (4) an exposure route (i.e., ingestion, inhalation, or dermal contact) at the exposure point (EPA, 1989a).

3.1.1 Contaminant Sources and Migration

Site investigations at the EKCO facility have confirmed the presence of VOCs in two water bearing units: the upper (shallow/intermediate) unit and the lower (bedrock) unit. Groundwater within the upper unit typically occurs under unconfined to semiconfined conditions; groundwater flow in this unit is toward wells W-1 and W-10. Groundwater in the lower unit is semi-confined by overlying shale; the direction of groundwater flow in this unit is also toward wells W-1 and W-10.

Chemicals detected in these groundwater units could migrate off site or persist on site. The migration of chemicals in groundwater is generally slower than groundwater migration itself, due to retardation of chemical movement resulting from adsorption to soils. This effect may be negligible for highly mobile chemicals (low adsorption potential) but could be extremely significant for the relatively immobile chemicals (high adsorption potential). In addition, chemicals disperse laterally as they are transported downgradient, resulting in dilution as contaminated water mixes with laterally adjacent, less contaminated water.

Chemicals in the groundwater could be transported into the air through volatilization or within the groundwater into off-site downgradient areas. Chemicals in groundwater can volatilize to the soil pore spaces and reach the surface of the soil and the surrounding air by diffusion. Volatilization from a water table aquifer is affected by depth to the water table and the moisture content of the unsaturated soil column. Volatilization from groundwater is most rapid if the water table is shallow. At the EKCO facility, the top of shallow groundwater unit averages about 15 feet below land surface. Volatilization to the surface from this aquifer is not expected because the distance to the land surface would greatly limit diffusion. In addition, rain and other precipitation events on the surface would result in water percolation downward through the soil. As water infiltrated through the soil, any chemicals in the air-filled soil pore spaces would become solubilized and be transported downward with the leachate back to the groundwater. However, if on-site groundwater were pumped to the surface for use, chemicals could volatilize directly to the air. Contaminated groundwater used as a domestic or nondomestic water supply could result in a significant loss of contaminants through volatilization to the atmosphere. Through agitation, heating, or other mechanical handling, VOCs would be expected to volatilize rapidly. Fate processes other than adsorption, dispersion, and volatilization (e.g., bioaccumulation, biodegradation, photodegradation) are expected to be insignificant in groundwater.

3.1.2 Receptors and Exposure Routes

The primary exposure pathway at the EKCO facility is the groundwater pathway. The EKCO facility is an operating manufacturing facility that produces nonstick bakeware. The

area surrounding the facility is largely urban and industrial, although land to the northwest is more rural with a large proportion of open space. Ohio Water Service (OWS) production wells are located 2,000 feet northeast of the EKCO property and are screened in the deep sand and gravel glacial valley. These wells provide water to the City of Massillon, Ohio.

Currently, groundwater on the EKCO property is used for production purposes at the facility. Workers at the site do not use (or plan to use in the future) this water for drinking or showering purposes. Presently, groundwater at the facility is pumped, air-stripped and then used for non-contact cooling. Therefore, workers are not exposed impacted groundwater, and were not evaluated in this risk assessment. Groundwater in the upper unit flows toward the production wells. Due to the operation of these wells, the shallow unit is dewatered beneath a large portion of the property. An intermediate unit exists only beneath the eastern portion of the property. A significant bedrock drawdown cone induced by the pumping of these production wells exists beneath the facility. Based on the current status of the EKCO facility, the only individuals expected to be exposed to site chemicals would be future on-site residents who use groundwater as their primary water supply.

In this risk assessment, on-site wells were evaluated for groundwater use by hypothetical future residents, assuming shutdown of the production wells and unrestricted future use of the property. Future residential exposure was evaluated for an adult and a small child (1-6 years) based on the fact that although the majority of a lifetime is spent as an adult, children are most sensitive to the effects of chemical exposure. Both shallow and bedrock well data were used to evaluate risks from drinking water ingestion, noningestion household uses (inhalation of and dermal contact with VOCs in tap water) and home garden watering uses (i.e., irrigation) by on-site future residents. It is noted that the upper unit has a very low-to-negligible probability as drinking water sources, since the yield from these units is not sufficient to support drinking water use (see Appendix C). Moreover, current off-site residences obtain drinking water from the bedrock or deep glacial sand and gravel aquifers. This issue is discussed in the Uncertainty Analysis (Section 6).

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3.2 **QUANTIFICATION OF EXPOSURE**

The degree of receptor exposure that occurs through each exposure pathway is determined by behavioral, chemical, and physiological factors. Behavioral factors affecting exposure would include the amount of time spent on site, the activities engaged in while on site, and the amount and type of clothing worn. Chemical factors affecting the degree of exposure would include the extent to which a chemical is absorbed through the skin (i.e., the absorption efficiency) or is volatilized to the air. Physiological factors affecting exposure would include the ability of the body to metabolize and eliminate the chemical(s). To quantify exposures in the risk assessment process, it is necessary to make assumptions concerning these factors in the absence of specific, detailed information. These assumptions are represented by a series of exposure parameters that quantify the magnitude, frequency, and duration of the exposure.

EPA has developed exposure algorithms for use in calculating chemical intakes (or doses) through specific exposure pathways and routes (EPA, 1989a). These algorithms combine the chemical exposure point concentrations with pathway- and route-specific parameters to produce daily chemical intakes in terms of the milligrams of chemical taken into the body per kilogram of body weight per day (mg/kg-day). For the EKCO facility, daily chemical intakes were calculated separately for carcinogenic and noncarcinogenic health effects under reasonable maximum exposure (RME) conditions, in accordance with EPA guidance (EPA, 1989a). The RME is the highest exposure that is reasonably expected to occur at a site.

The following paragraphs discuss the exposure algorithms that were used to calculate chemical intakes in this study.

3.2.1 Groundwater Ingestion

Daily chemical intakes through groundwater ingestion for adults and children were calculated using the equation shown in Table 3-1. The values for the groundwater ingestion exposure parameters that were used are also shown in Table 3-1.

The ingestion rate for potable water is established at 2 L/day for an adult (EPA, 1991a) and 1 L/day for a child (EPA, 1989a; 1989b) for a residential land use. These amounts include water consumed in the form of other beverages and the ingestion of foods prepared in or with water.

The exposure frequency (EF) for the potential resident (i.e., 350 days/year), is the recommended residential exposure frequency (EPA, 1991a). This value assumes that 15 days per year are spent away from home.

The exposure duration (ED) for an adult resident was assumed to be a maximum of 24 years. The ED for the child resident was assumed to be a maximum of 6 years. These values are based on the 90th percentile total number of years (i.e., 30) at one residence (EPA, 1989a), assuming conservatively that the initial 6 years of exposure occur during early childhood when the sensitivity to chemical exposure is the greatest.

The value for body weight is the average body weight over the exposure period. An average body weight is used because when combined with other variable values in the intake equation, it is believed to result in the reasonable maximum exposure. Incorporating a higher body weight with the same intake rate would result in lower exposure than the reasonable maximum. In addition, using an average body weight rather than a reasonable maximum is recommended because the available toxicity values are based on average body weight. The recommended average body weight for an 18- to 75-year old adult is 70 kg; the recommended average body weight for a 1- to 6-year-old child is 15 kg (EPA, 1991a).

TABLE 3-1

EXPOSURE ALGORITHM AND ASSUMPTIONS FOR POTENTIAL INGESTION OF CHEMICALS IN GROUNDWATER EKCO HOUSEWARES MASSILLON, OHIO

EXPOSURE ALGORITHM

Intake from Water Ingestion (mg/kg-day) = $\frac{CW \times IR \times CF \times EF \times ED}{BW \times AT}$

where:

CW = Chemical concentration in water $(\mu g/L)^a$

IR = Water ingestion rate (L/day) CF = Conversion factor (10^{-3} mg/ μ g) EF = Exposure frequency (days/year)

ED = Exposure duration (years)

BW = Body weight (kg) AT = Averaging time (days)

EXPOSURE ASSUMPTIONS b

	Receptor Group Resident				
Variable	Child	Adult			
IR (L/day)	1	2			
CF (mg/μg)	1E-3	1E-3			
EF (days/year)	350	350			
ED (years) ^c	6	24			
BW (kg)	15	70			
AT-Noncancer (days)	2,190	8,760			
AT-Cancer (days)	25,550	25,550			

^a See Tables 2-1 and 2-2 for chemical-specific exposure point concentrations.

^b Default exposure assumptions from U.S. EPA (1991a), unless indicated.

^c Based on 95th percentile time at one residence (U.S. EPA, 1989a), assuming 20% of time is as a child (6 years out of 30 years).

The averaging time (AT) is the period over which the exposure is averaged. For noncarcinogenic effects, AT is equal to the exposure duration (ED). For carcinogenic effects, AT is equal to the average lifetime for an individual in the U.S., which was assumed to be 70 years (EPA, 1991a).

The estimated daily chemical intakes that result from groundwater ingestion are presented in Appendix B, Tables B-1 and B-5.

3.2.2 Household Noningestion Uses of Groundwater (Inhalation of VOCs)

Household water is used for noningestion purposes such as showering, dish washing, laundering, cooking, and bathing. Volatile contaminants in the water may enter the indoor air through these uses, exposing receptors to volatilized contaminants through inhalation. Table 3-2 presents the equation and exposure parameters used to calculate daily chemical doses via inhalation of volatiles from groundwater (upper and lower units). This equation contains a volatilization constant (K) which has an upper bound value of 5E-01 L/m³. The volatilization constant was derived by Andelman (as presented in EPA, 1991b). It defines the relationship between the concentration of a contaminant in household water and the average concentration of the volatilized contaminant in household air. The constant is based on the following assumptions:

- volume of water used by a hypothetical family of four is 720 L/day;
- volume of the hypothetical residential dwelling is 150,000 L;
- air exchange rate is 0.25 m³/hr;
- an average of 50% of the chemical concentration in groundwater is transferred into household air from all water uses.

According to EPA (1991b), VOCs are defined as organic chemicals with a molecular weight of less than 200 g/mol, and a Henry's Law Constant of greater than 1E-05 atm-m³/mol.

TABLE 3-2

EXPOSURE ALGORITHM AND ASSUMPTIONS FOR POTENTIAL NONINGESTION (INHALATION) HOUSEHOLD USES OF GROUNDWATER EKCO HOUSEWARES MASSILLON, OHIO

EXPOSURE ALGORITHM

Inhalation from Household Water Use	_	$CW \times K \times IR \times EF \times ED$
(mg/k-d)	_	$\overline{BW \times AT}$

where:

CW = Chemical concentration in water $(\mu g/L)^a$

K = Volatilization Factor (L/m³)

IR = Daily indoor inhalation rate (m^3/day)

CF = Conversion factor $(10^3 \text{ mg/}\mu\text{g})$ EF = Exposure frequency (days/year)

ED = Exposure duration (years)

BW = Body weight (kg)

AT = Averaging time (days)

EXPOSURE ASSUMPTIONS^b

	Receptor Group				
	Resi	dent			
Variable	Child	Adult			
K L/m³)°	5E-1	5 E -1			
IR (m³/day) ^c CF (mg/μg)	15 1E-3	15 1E-3			
EF (days/year)	350	350			
ED (years) ^c	6	24			
BW (kg)	15	70			
AT-Noncancer (days)	2,190	8,760			
AT-Cancer (days)	25,550	25,550			

^a See Tables 2-1 and 2-2 for chemical-specific exposure point concentrations for upper and lower units, respectively.

^b Default exposure assumptions from U.S. EPA (1991a), unless indicated.

^c Default values based on inhalation equation derived for household water use (EPA, 1991b). The constant K is actually unitless, but incorporate a conversion factor to estimate air concentration in mg/m³.

^d Based on 95th percentile time at one residence (U.S. EPA, 1989a), assuming 20% of time is as a child (6 years out of 30 years).

An inhalation rate (IR) of 15 m³/day has been found to represent a reasonable upperbound inhalation rate for daily indoor residential activities (EPA, 1991a). The values used for the remaining exposure parameters in the inhalation equation (i.e., EF, ED, BW, and AT) were discussed in Subsection 3.2.1.

The estimated daily chemical intakes that result from inhalation of volatiles are presented in Appendix B, Tables B-3 and B-7.

3.2.3 Dermal Absorption of VOCs While Showering/Bathing

For most contaminants, dermal contact with water during bathing or swimming will generally pose less threat than direct consumption of water, since contact times are relatively brief and most chemicals do not readily penetrate the skin when present at dilute levels in water (EPA, 1989b). In addition, VOCs have been shown to volatilize significantly during showering (Foster and Chrostowski, 1987). These observations suggest that the effective concentrations of VOCs in water available for dermal absorption will be relatively small (EPA, 1992). Nevertheless, the potential exposure by this pathway was estimated for future residential use.

Table 3-3 presents the equation used to calculate daily chemical intake via dermal absorption while bathing. The exposure input parameters used are detailed below the table. The surface area for the adult represents 100% of the maximum surface area (23,000 cm²) for the male (Table 8-3; EPA, 1992). For the child, the 95th percentile value for total surface area was used (8,760 cm²) (Table 8-4; EPA 1992). It was assumed that both the child and adult shower once daily for 15 minutes giving an exposure time (ET) of 0.25 hours per day for 350 days per year (EPA, 1992). The permeability constant (Kp) defines the rate of movement of a waterborne chemical through the skin. Chemical-specific permeability constants for the chemicals of concern were obtained from EPA (1992), and are presented in Table 3-4.

Table 3-3

Exposure Algorithm and Assumptions for Potential Dermal Absorption While Showering/Bathing Ekco Housewares Massillon, Ohio

Dermal Absorption Dose from Surface Water	$= \frac{\mathbf{CW} \times \mathbf{CF} \times \mathbf{SA} \times \mathbf{Kp} \times \mathbf{ET} \times \mathbf{EF} \times \mathbf{ED}}{\mathbf{EF} \times \mathbf{ED}}$
(mg/kg-day)	BW×AT

Where:

CW = Chemical concentration in surface water (mg/L)

CF = Conversion factor (10⁻³ L/cm³)

SA = Skin surface area available for contact (cm²/day)

Kp = Dermal permeability constant (cm/hour)

ET = Exposure time (hours/day)

EF = Exposure frequency (days/year)

ED = Exposure duration (years)

BW = Body weight (kg)

AT = Averaging time (days)

Exposure Assumptions (child and adult on-site resident):

CW = 95% UCL in upper and lower groundwater units.

 $SA = 8,760 \text{ cm}^2/\text{day for the child resident (EPA, 1994)}.$

= 23,000 cm²/day for the adult resident (EPA, 1994).

Kp = Chemical specific (see Table 3-4).

ET = 0.25 hr/day (EPA, 1994).

EF = 350 days/year (EPA, 1989).

ED = 6 yrs for the child resident (EPA, 1991).

= 24 yrs for the adult resident (EPA, 1991).

BW = 15 kg for the child (EPA, 1991).

= 70 kg for the adult (EPA, 1991).

AT = Exposure duration (years) x 365 days/year for evaluating noncancer risk.

= 70 years x 365 days/year for evaluating cancer risk.

The values used for the remaining exposure parameters in the dermal absorption equation (i.e., EF, ED, BW, and AT) were discussed in Subsection 3.2.1.

The estimated daily chemical doses that result from dermal absorption from groundwater are presented in Appendix B, Tables B-2 and B-6.

3.2.4 Consumption of Home-Grown Produce

Exposure to site contaminants through the ingestion of garden vegetables and fruits irrigated with contaminated groundwater is a potential exposure route for future child and adult residents. This potential exposure route was evaluated for the shallow and bedrock groundwater units. Three general categories of garden produce were considered in this evaluation -- leafy vegetables (represented by lettuce), root vegetables (represented by carrots), and garden fruits (represented by tomatoes). The equation and assumptions used to estimate chemical intakes through the consumption of home-grown produce are presented in Table 3-5.

EPA (1991a) has recommended intake rates of 42 g/day for consumption of home-grown fruit and 80 g/day for adult consumption of home-grown vegetables. These values have been adjusted to reflect the amount of home-grown produce typically consumed as a proportion of total consumption (i.e, 30% for fruits and 40% for vegetables). The total consumption of fruits and vegetables by a child (1-6 years) was assumed to be 60% of the total consumption of fruits and vegetables by an adult. This assumption is based on data collected by USDA on the national average consumption rates per day for lettuce, tomatoes, and carrots (Pao et al., 1982).

For an adult, it was assumed that leafy vegetables and root vegetables comprise 45% and 55% of the total vegetable consumption, respectively. For a child, it was assumed that leafy vegetables and root vegetables comprise 35% and 65% of the total vegetable consumption, respectively. These assumptions were based on tomato, carrot, and lettuce consumption data for 19- to 34-year-old (male) and 3- to 5-year-old (male and female) individuals (Pao

Table 3-4

Dermal Permeability Coefficients for Chemicals of Concern at EKCO Housewares

Chemical	Groundwater Unit	Kp* (cm/hr)
Acetone	S	NA
Benzene	S	2.1E-02
Carbon disulfide	S	2.4E-02
Chloroethane	S	8.0E-03
Chloroform	S	8.9E-03
1,1-Dichloroethane	S,B	8.9E-03
1,2-Dichloroethane	S	5.3E-03
1,1-Dichloroethene	S,B	1.6E-02
1,2-Dichloroethene	S,B	1.0E-02
Ethylbenzene	S	7.4E-02
Methyl ethyl ketone	S	1.1E-03
4-Methyl-2-pentanone	S	3.3E-03
Tetrachloroethene	S	4.8E-02
Toluene	S	4.5E-02
1,1,1-Trichloroethane	S,B	1.7E-02
1,1,2-Trichloroethane	S	8.4E-03
Trichloroethene	S,B	1.6E-02
Vinyl chloride	S,B	7.3E-03
Xylene (m)	S	8.0E-02

NA = Not available

S = present in shallow unit

B = present in bedrock unit

*Kp value obtained from Table 5-7 (EPA, 1992)

TABLE 3-5

EXPOSURE ALGORITHM AND ASSUMPTIONS FOR POTENTIAL INGESTION OF CHEMICALS IN HOME-GROWN PRODUCE EKCO HOUSEWARES MASSILLON, OHIO

EXPOSURE ALGORITHM

Intake from Home-GrownProduce $\frac{[(C_{lv} \times IR_{lv}) + (C_{gf} \times IR_{gf}) + (C_{rv} \times IR_{rv})] \times EF \times ED}{BW \times AT}$

where:

 C_{lv} = Chemical concentration in leafy vegetables $(mg/kg)^a$

 IR_{lv} = Ingestion rate for leafy vegetables (kg/day)

C_{gf} = Chemical concentration in garden fruits (mg/kg)^a

 IR_{ef} = Ingestion rate for garden fruits (kg/day)

 C_{rv} = Chemical concentration in root vegetables $(mg/kg)^a$

 IR_{rv} = Ingestion rate for root vegetables (kg/day)

EF = Exposure frequency (days/year)

ED = Exposure duration (years)

BW = Body weight (kg) AT = Averaging time (days)

EXPOSURE ASSUMPTIONS^b

	Receptor Group Resident				
V ariable	Child	Adult			
IR _{Iv} (kg/day) ^c	0.017	0.036			
IR _{gf} (kg/day) ^c	0.025	0.042			
IR _{rv} (kg/day) ^c	0.031	0.044			
EF (days/year)	350	350			
ED (years) ^d	6	24			
BW (kg)	15	70			
AT-Noncancer (days)	2,190	8,760			
AT-Cancer (days)	25,550	25,550			

^a See Table 3-4 for calculation of chemical concentrations in fruits and vegetables.

^b Default exposure assumptions from U.S. EPA (1991b), unless indicated.

[&]quot;Based on U.S. EPA (1991b) and Pao, et al (1982).

^d Based on 95th percententile time at one residence (U.S. EPA, 1989a), assuming 20% of time is as a child (6 years out of 30 years).

The values used for the remaining exposure parameters in the produce ingestion equation (i.e., EF, ED, BW, and AT) were discussed in Subsection 3.2.1.

The estimated daily chemical intakes that result from ingestion of garden produce are presented in Appendix B.

Concentrations of chemicals in garden produce were not measured. Rather, chemical concentrations in garden fruits and vegetables were estimated based on the assumption that the plants accumulate chemicals from soil that has been irrigated with contaminated groundwater from the shallow or bedrock units.

The chemical concentrations in the edible plant portions were calculated as follows:

where:

CS = Chemical concentration in soil (mg/kg). PUF = Plant uptake factor (unitless).

The plant uptake factor relates the chemical concentration in the plant to the chemical concentration in the soil under equilibrium conditions. Although plants take up chemicals from the soil solution, uptake factors are generally derived for soil solids. To estimate the chemical concentration in soil resulting from waterings of gardens, the shallow groundwater chemical concentration was multiplied by the chemical-specific soil-water partition coefficient (K_d). The soil-water partition coefficient relates the chemical concentration adsorbed to the solid phase to the chemical concentration in the soil solution. If a K_d value was not available for a chemical, it was estimated from the organic-carbon partition coefficient (K_∞) and the soil organic carbon content (f_∞). The concentrations of VOCs in groundwater were assumed to remain constant over time; loss of VOCs due to volatilization to the atmosphere was not considered in estimating soil concentrations.

Two approaches were used to determine plant uptake factors. One approach was used for leafy vegetables and garden fruits, and the other was used for root vegetables. These approaches are discussed below.

3.2.3.1 Uptake Factors for Leafy Vegetables and Garden Fruits

The uptake factors for VOCs in leafy vegetables and garden fruits were derived from the octanol-water partition coefficient for each chemical, applying the following relationship from Travis and Arms (1988):

$$\log B = 1.588 - 0.578 \log K_{ow}$$

where:

B = Bioconcentration factor for vegetation: the ratio of the chemical concentration in above-ground plant parts (mg chemical per kg of dry plant) to the chemical concentration in soil (mg chemical per kg of dry soil).

 $Log K_{ow} = Log of the octanol-water partition coefficient.$

The uptake factor was then calculated by multiplying B by the percent dry weight of the plant:

$$PUF = B \times \%$$
 Dry Weight

Since vegetable/fruit ingestion rates are expressed in terms of wet plant weight, chemical concentrations calculated using the Travis and Arms equation must be converted from a dry weight to a wet weight basis using the plant moisture content. Leafy vegetables and garden fruits typically have a water content of 95 to 96 percent (Baes et al., 1984). A water content of 95 percent was used in this analysis, which is equivalent to 5 percent dry weight (0.05). It is important to note that the Travis and Arms equation is based on data for 29 chemicals, each with a $\log K_{ow}$ value falling in the range from 1.15 to 6.89; the use of this equation may be inappropriate for chemicals with values that fall outside this range.

3.2.3.2 Uptake Factors for Root Vegetables

Uptake factors for root vegetables were derived based on the work of Briggs et al. (1982). Briggs studied the uptake of organic chemicals from solution by barley shoots and established the following relationship between the root concentration factor (RCF) and the K_{ow} for the organics tested:

$$log (RCF - 0.82) = 0.77 log K_{ow} - 1.52$$

where:

 $RCF = C_{root}/C_{soln}$

 C_{root} = Chemical concentration in the root (mg/kg). C_{soln} = Chemical concentration in solution (mg/L).

The chemical distribution in the soil and water phases can be described as:

$$C_{\text{soil}}/C_{\text{soln}} = K_{\infty} \times f_{\infty}$$

The plant uptake factor for root vegetables (RUF) for each compound can be determined from the RCF, given the following relationship:

$$RUF = \frac{C_{root}/C_{soln}}{C_{soil}/C_{soln}} = \frac{C_{root}}{C_{soil}} = \frac{RCF}{K_{\infty} \times f_{\infty}}$$

where:

 C_{root} = Concentration in root (mg/kg).

 C_{soil} = Concentration in soil (mg/kg).

 C_{soln} = Concentration in solution (mg/L).

 K_{∞} = Organic carbon partition coefficient.

 f_{∞} = Fraction of organic carbon in soil.

RCF = Root concentration factor, or the ratio of the root concentration to the solution concentration.

The estimated chemical concentrations in leafy vegetables, garden fruits, and root vegetables are presented in Tables 3-6 and 3-7, along with the chemical-specific parameters used in this derivation.

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Table 3-6

Estimation of Chemical Concentrations in Garden Produce
Upper Unit

		Octanol-Water	Organic Carbon	Soil-Water	Root Vegetables		Leafy V	Leafy Vegetables/Garden Fruits		
Chemical	Groundwater Concentration (µg/L)	Partition Coefficient (log K _{ow})	Partition Coefficient (log K _∞)	Partition Coefficient (K _d)	RCF (L/kg)	Concentration (mg/kg)	В	PUF	Concentration (mg/kg)	
1,1-Dichloroethane	451	1.78	1.48	0.60	1.53	6.90E-01	3.62	0.18	4.94E-02	
1,1-Dichloroethene	220	1.92	1.81	1.29	1.73	3.80E-01	3.01	0.15	4.27E-02	
1,2-Dichloroethene (total)	93.7	2.09	1.77	1.18	2.05	1.92E-01	2.40	0.12	1.32E-02	
1,1,1-Trichloroethane	5380	2.38	2.11	2.58	2.87	1.55E+01	1.63	0.08	1.13E+00	
Trichloroethene	22300	2.89	1.99	1.95	5.89	1.31E+02	0.83	0.04	1.80E+00	
Vinyl chloride	28.4	0.6	0.39	0.05	0.91	2.58E-02	17.43	0.87	1.21E-03	
Chloroethane	3.19	1.43	0.51	0.06	1.20	3.83E-03	5.77	0.29	5.96E-05	
Acetone	4.25	-0.24	-0.43	0.01	0.84	3.57已-03	53.30	2.66	8.42E-05	
1,2-Dichloroethane	9.24	1.46	1.22	0.33	1.22	1.13E-02	5.55	0.28	8.51E-04	
2-Butanone	2.6	0.28	0.09	0.02	0.87	2.26E-04	26.68	1.33	8.53E-05	
1,1,2-Trichloroethane	15.61	2.18	1.75	1.12	2.26	3.53E-02	2.13	0.11	1.87E-03	
Tetrachioroethene	7.71	2.61	2.44	5.51	3.91	3.01E-02	1.20	0.06	2.55E-03	
Toluene	12.6	2.62	2.12	2.64	3.96	4.99E-01	1.18	0.06	1.97E-03	
Ethylbenzene	1.41	3.08	2.25	3.56	7.93	1.12E-03	0.64	0.03	1.61E-05	

Table 3-6
Estimation of Chemical Concentrations in Garden Produce
Upper Unit

	Octanol-Water		Organic Carbon Soil-Water	Root Vegetables		Leafy Vegetables/Garden Fruits			
Chemical	$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	Coefficient	RCF (L/kg)	Concentration (mg/kg)	В	PUF	Concentration (mg/kg)		
Benzene	1.45	2.04	1.91	1.63	1.94	2.82E-03	2.56	0.13	3.02E-04
Xylenes (total)	1.56	3.1	2.81	12.91	8.18	1.28E-03	0.63	0.03	6.30E-04
Carbon disulfide	1.1	2.03	2.47	5.90	1.92	2.12E-04	2.60	0.13	8.43E-04
Chloroform	1.99	1.94	1.64	0.87	1.76	3.51E-03	2.93	0.15	2.54E-04
4-Methyl-2-pentanone	2.92	1.09	0.79	0.12	1.03	3.00E-03	9.08	0.45	1.63E-04

^{*} Exposure point concentrations for upper unit (See Table 2-1)

 $K_d = K_{oc} \times f_{oc}$ (f_{oc} assumed to be 0.02) log (RCF - 0.82) = (0.77 log K_{ow}) - 1.52 (Briggs, et al, 1982) log B = 1.588 - (0.578 log K_{ow}) (Travis and Arms, 1988) PUF = B x % dry weight

Table 3-7

Estimation of Chemical Concentrations in Garden Produce Lower Unit EKCO Housewares Massillon, Ohio

Chemical	Octanol-Water		Organic Carbon Soil-Water		Root Vegetables		Leafy Vegetables/Garden Fruits		
	Concentration Coeffici	Partition Coefficient ^b (log K _{ow})	Partition Coefficient ^b (log K _∞)	Partition Coefficient (K _d)	RCF (L/kg)	Concentration (mg/kg)	В	PUF	Concentration (mg/kg)
1,1-Dichloroethane	102	1.78	1.48	0.60	1.53	1.55E-01	3.62	0.18	1.11E-02
1,1-Dichloroethene	31.3	1.92	1.81	1.29	1.73	5.40E-02	3.01	0.15	6.07E-03
1,2-Dichloroethene (total)	142	2.09	1.77	1.18	2.05	2.92E-01	2.40	0.12	2.01E-02
1,1,1-Trichloroethane	613	2.38	2.11	2.58	2.87	1.76E+00	1.63	0.08	1.29E-01
Trichloroethene	234	2.89	1.99	1.95	5.89	1.38E+00	0.83	0.04	1.89E-02
Vinyl chloride	8.15	0.6	0.39	0.05	0.91	7.40E-03	17.43	0.87	3.49E-04

^{*} Exposure point concentrations for lower unit (see Table 2-2).

Calculation Notes:

 $K_d = K_{\infty} \times f_{\infty} (f_{\infty} \text{ assumed to be } 0.02)$

 $log (RCF - 0.82) = (0.77 log K_{ow}) - 1.52 (Briggs, et al, 1982)$

 $\log B = 1.588 - (0.578 \log K_{ow})$ (Travis and Arms, 1988)

PUF = B x % dry weight

^b Values from Montgomery and Welkom, 1991.

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SECTION 4 TOXICITY ASSESSMENT

4.1 <u>INTRODUCTION</u>

The purpose of this section of the report is to explain how the toxicity criteria for the chemicals evaluated for the EKCO site are derived and used in this baseline risk assessment. Toxicity criteria are applied to the estimated doses for the potentially exposed human calculated in Section 3 to determine potential adverse health effects. Human health criteria (cancer slope factors and reference doses) are developed and approved for most chemicals by the EPA, and are described in the Integrated Risk Information System (IRIS), an EPA computerized database (IRIS, 1994). If not available through IRIS, data was obtained from the Health Effects Assessment Summary Tables (HEAST; EPA, 1994). For some chemicals not listed in IRIS or HEAST, the criteria were obtained through the Environmental Criteria Assessment Office (ECAO), as recommended by EPA (1989) in their risk assessment guidance.

In predicting potential human health risks, both cancer (carcinogenic) and noncancer (noncarcinogenic) health effects of a chemical must be considered. The potential for producing carcinogenic effects is limited to those substances shown to cause cancer in animals and/or humans. Excessive exposure to all substances, carcinogens and noncarcinogens, can produce adverse noncarcinogenic effects. Therefore, cancer slope factors (CSFs) are developed for those chemicals of concern classified as carcinogens, and reference doses (RfDs) are developed to express the noncancer effects for all chemicals of concern selected regardless of its carcinogenic classification.

Animal data are the most common source of toxicity information although human response data are preferred. In general, much higher doses of a chemical in animals are required to produce an equal toxic response in humans. General principles of extrapolation for carcinogens and noncarcinogens can be found in later subsections of this toxicity assessment.

The toxicity of a chemical depends on the dose, route of administration, duration and frequency of exposure, and the species evaluated. The term "toxicity" is a property of the chemical that causes varying degrees of morphological and/or biochemical tissue or organ damage as the dose increases. The lower a dose of chemical that produces a given adverse effect, the more toxic or potent the chemical is considered. Some chemicals may produce toxic effects ranging from respiratory/skin irritation to lethality after a single (acute) high dose. However, acute exposures are not usually the main focus of concern in the baseline risk assessment. Exposure for a continuous period of months or years (chronic) at low exposure levels are potentially more significant from a human health viewpoint.

The probability of occurrence and the severity of a chronic toxic effect also depend on the concentration of chemical present in various media (water, soil, air) as well as the amount that can be accumulated by a person (exposure dose) on or near the site. Chemicals are potentially capable of producing adverse effects through breathing (inhalation), ingestion by mouth (oral ingestion), or skin absorption (dermal contact). Some chemicals may produce toxicity only through one route. Others may cause toxicity through a combination of some or all routes. Consequently, each chemical is evaluated for toxicity by determining its potency through each exposure route.

Last of all, a given species (e.g., mouse, rat, dog, human) may exhibit different types or degrees of toxicity to a given dose of a chemical. These qualitative and quantitative species differences in toxic response are related to genetic and other physiological factors such as rates and amounts of:

- absorption by different routes.
- distribution to various organs and tissues.
- metabolism of the chemical by the liver.
- excretion by the kidneys or lungs.

Moreover, there may be interspecies differences in chemical sensitivity, i.e., the magnitude of damage to a given tissue or organ produced by the same amount of chemical may vary

in different species. Finally, some chemicals may affect entirely different organ systems in different species.

4.2 CARCINOGENS

4.2.1 Classification of Chemicals Inducing Cancer - Weight of Evidence

Carcinogens are evaluated in a two-phased approach by the EPA (Carcinogenic Risk Assessment Verification Endeavor; CRAVE). First, their weight of evidence for causing cancer is determined, and then their CSF is determined if sufficient data are available. Both human and animal cancer data are reviewed to determine the likelihood that a chemical is a human and/or animal carcinogen. These weight of evidence classifications are summarized in Table 4-1. Specific EPA classifications for the cancer-causing contaminants of concern detected at the EKCO site are presented in Table 4-2.

4.2.2 Potency of Cancer-Causing Chemicals

Following weight of evidence classification, the potency of the chemical at which cancer is thought to occur in humans is calculated. The carcinogenic potency of a chemical is defined as that dose producing cancer over a lifetime of exposure by a given exposure route. A CSF is used to express this potency.

CSFs are expressed either as risk per unit dose ([mg/kg-day]-¹) or risk per unit concentration ([mg/L or mg/m³]-¹). They represent values that quantitatively define the relationship between exposure dose and carcinogenic response for a given chemical. By definition, the CSF represents a plausible upper-bound estimate of the probability of a response per unit intake of a chemical during an entire average lifetime of exposure (70 years). The larger the CSF for a given carcinogen, the greater is the risk of cancer occurring at a specific exposure level. CSFs may be developed from controlled studies in laboratory animals or epidemiologic studies in humans. The most valid data set is used to determine the CSF. If animal data are used, a species is chosen that has the closest physiological response to humans. When no choice is obvious, the most sensitive species, or an average of several supportive data sets, is used.

Table 4-1

EPA Categorization of Carcinogens
Based on Human and Animal Evidence

EPA Categorization of Carcinogens (EPA, 1986b)										
Animal Evidence										
	Sufficient	Limited	Inadequate	No Data	No Evidence					
Human Evidence		,								
Sufficient	Α	Α	Α	Α	Α					
Limited	В1	В1	B1	B 1	В1					
Inadequate	B2	С	D	D	D					
No data	В2	С	D	D	E					
No evidence	B2	С	D	D	E					

Key:

- Group A Human carcinogen (sufficient evidence from epidemiological studies).
- Group B1 Probable human carcinogen (at least limited evidence of carcinogenicity to humans).
- Group B2 Probable human carcinogen (a combination of sufficient evidence in animals and inadequate data in humans).
- Group C Possible human carcinogen (limited evidence in animals in the absence of human data).
- Group D Not classified (inadequate animal and human data).
- Group E No evidence for carcinogenicity (no evidence for carcinogenicity in at least two adequate animal tests in different species, or in both epidemiological and animal studies).

Table 4-2

Oral and Inhalation Cancer Slope Factors EKCO Housewares Massillon, Ohio

Chemical	Weight-of- Evidence Classification	Oral Slope Factor (mg/kg-day)	Species/ Type of Cancer	Stope Factor Basis/ Source	Inhalation Slope Factor (mg/kg-day)	Species/ Type of Cancer	Slope Factor Basis/ Source
Benzene	А	2.9E-02	human/leukemia	occupational exposure/IRIS	2.9E-02	human/leukemia	oral slope factor/IIEAST
Chloroform	B2	6.1E-03	rat/kidney	104-week drinking water study/IRIS	8.1E-02	mouse/liver	gavage/HEAST
1,1-Dichloroethane	С	ND		/IRIS	ND		/IRIS
1,2-Dichloroethane	B2	9.1E-02	rat/circulatory system	gavage/IRIS	9.1E-02	rat/circulatory system	oral slope factor/HEAST
1,1-Dichloroethene	С	6.0E-01	rat/adrenal	gavage/IRIS	1.2E+00	mouse/kidney	12-month inhalation study/HEAST
Tetrachloroethene	B2/C	5.2E-02	NA	NA/ECAO	2E-03	NA	NA/ECAO
1,1,2-Trichloroethane	С	5.7E-02	mouse/liver	gavage/IRIS	5.7E-02	mouse/liver	oral slope factor/HEAST
Trichloroethene	B2/C	1.1E-02	NA	NA/ECAO	6E-03	NA	NA/ECAO
Vinyl Chloride	Α	1.9E+00	rat/lung,liver	1000-day dietary study/HEAST	3.0E-01	rat/liver	1-year inhalation study/ΗΕΑSΤ

IRIS - Integrated Risk Information System
HEAST - Health Effects Assessment Summary Tables ECAO - Environmental Criteria and Assessment Office

ND - Not determined NA - Not available

In animal studies designed to evaluate cancer-causing potential, relatively high doses of the chemical are usually administered daily for the lifetime of the animal (e.g., rat = 2 years). These daily doses are usually close to those producing noncarcinogenic toxicity. Cancer formation is more likely to be induced at these high doses in small groups of animals. Therefore, cancer effects occurring at higher animal doses must be extrapolated downward and consequently, the risk determined for a given carcinogen is likely to be a plausible upper limit. The actual (absolute) risk of cancer is unknown, but is likely to be considerably lower than the predicted (relative) risk, and may even be as low as zero (EPA, 1989). Moreover, in contrast to noncarcinogens, where it is believed that a safe dose exists below a threshold dose for toxicity, it is assumed that any dose of a carcinogen (however low) may cause changes in a single cell that could result in uncontrolled cell division, eventually leading to cancer. It has been argued that at low doses, cells may have the ability to detoxify carcinogens or repair chemical-induced cellular damage. Although it is important to recognize the possibility that some carcinogens may have a threshold for cancer-causing effects, it is assumed that no threshold exists when estimating cancer risk for humans. That is, it is conservatively assumed there is no dose of a carcinogen that is risk free. Figure 4-1 illustrates the "no-threshold" concept for carcinogens.

4.2.3 Route of Entry Into The Body

The carcinogenic potency of a substance depends on its route of entry into the body (e.g., ingestion, inhalation or dermal contact) and CSFs are, therefore, developed accordingly. In some cases, a carcinogen may produce tumors only at or near a specific route of entry (e.g., nasal passages) and may not be carcinogenic through other exposure routes. Table 4-2 presents the CSFs by oral ingestion and inhalation for the carcinogens of concern at the EKCO facility. Dermal CSFs are presented in Table 4-4.

4.2.3.1 Oral Ingestion Route

Oral slope factors are used to evaluate the risk from exposure to potential carcinogens through oral exposure pathways such as the incidental ingestion of soil and sediment, and the ingestion of groundwater (drinking water).

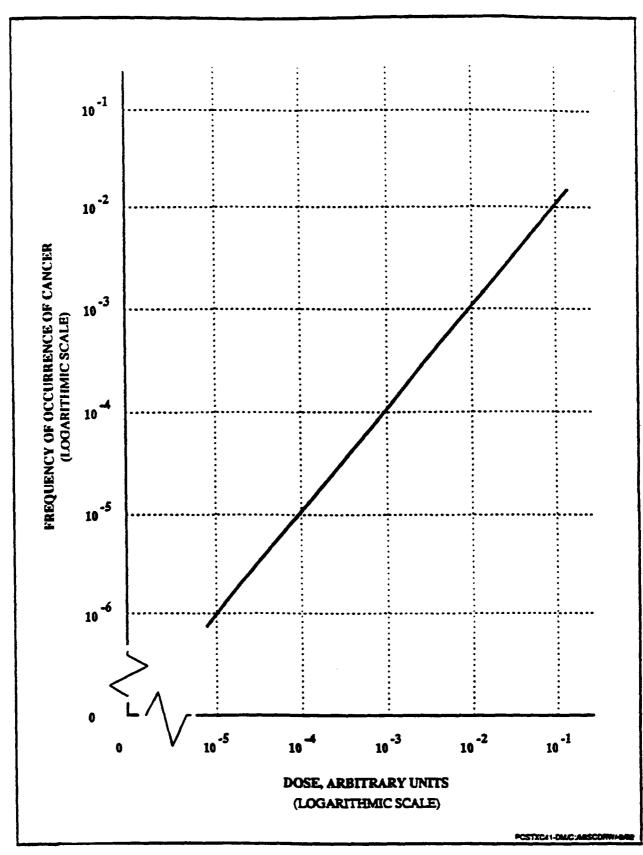


FIGURE 4-1 HYPOTHETICAL DOSE-RESPONSE CURVE FOR A "NO THRESHOLD" OR CARCINOGENIC CHEMICAL

4.2.3.2 Inhalation Route

Inhalation slope factors are used to evaluate the risk from exposure to potential carcinogens through pathways such as the inhalation of volatile chemicals from groundwater while showering. In the absence of an inhalation slope factor, oral slope factors, if available for a chemical, were used.

4.2.3.3 Dermal Absorption Route

Dermal slope factors are not available from the EPA, but it is assumed that chemicals which are carcinogenic orally could potentially produce cancer through dermal exposure. In the absence of dermal slope factors, the oral slope factor is divided by an appropriate gastrointestinal (GI) absorption factor (EPA, 1989). The GI absorption factor adjusts the orally administered dose upward for the amount absorbed since dermal exposure doses are expressed as "absorbed" doses. Oral and inhalation doses are usually expressed as "administered" doses.

Oral slope factors are normally developed from long-term studies where a substance is administered orally by gavage or diet to laboratory animals. Depending on the method and form in which the chemical is administered, the relative absorption of the chemical through the gastrointestinal tract (and therefore the relative absorption factor) may vary considerably. A conservative absorption factor of 80 percent for volatile organics was used. Metals, in general, tend to be poorly absorbed through the GI tract and therefore have the lowest absorption factor. The oral slope factor was divided by 0.8 (volatiles) to obtain the dermal slope factors.

4.3 **NONCARCINOGENS**

4.3.1 Estimates of Noncarcinogenic Potency and Toxicity

Toxicity criteria used to evaluate potential noncarcinogenic health effects are termed reference doses (RfDs). Unlike the approach used in evaluating carcinogenic risk, it is

assumed for noncarcinogenic chemicals that a threshold dose exists below which there is no potential for human toxicity.

The concept of the RfD was developed by the EPA to refer to the daily intake of a chemical to which an individual can be exposed without any expectation of noncarcinogenic effects (e.g., organ damage, biochemical alterations, birth defects) occurring during a given exposure duration.

The RfD is derived from a no-observed-adverse-effect level (NOAEL) or lowest-observed-adverse-effect level (LOAEL) obtained from human or animal studies. The criteria for choosing the appropriate NOAEL or LOAEL are discussed in the U.S. EPA's risk assessment guidance (EPA, 1989) and other documents (EPA, 1988). Rather than linearly extrapolating from high dose to low dose to obtain toxicity criteria as with carcinogens, RfDs are derived by applying standard uncertainty factors to the NOAEL or LOAEL, and in some cases, an additional modifying factor to account for professional assessment of scientific uncertainties in the available data (EPA, 1988).

A NOAEL is the highest dose of chemical at which <u>no</u> toxic effects are observed in any of the test subjects or animals. The study chosen to establish the NOAEL is based on the criterion that the measured toxic endpoint represents the most sensitive ("critical") target organ or tissue to the toxicity of that chemical (i.e., that target organ or tissue that shows evidence of damage at the lowest dose).

In contrast to a NOAEL, a LOAEL is the lowest dose at which any possible toxic effect is observed in any of the test subjects or animals. LOAELs are used to derive an RfD in the absence of a suitable NOAEL. Since it is possible that toxicity could occur at a lower dose than the reported LOAEL, an additional uncertainty factor is applied. Figure 4-2 illustrates the toxicity threshold concept as applied to a dose-response curve for a noncarcinogenic chemical.

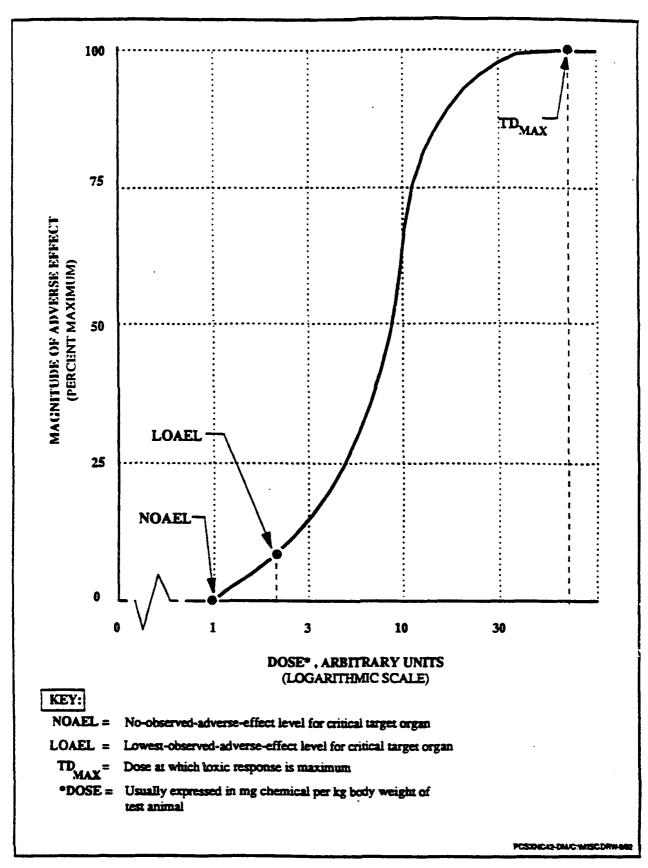


FIGURE 4-2 HYPOTHETICAL DOSE-RESPONSE CURVE FOR A "THRESHOLD" OR NONCARCINOGENIC CHEMICAL

4.3.2 Reference Doses

As with carcinogens, the route of exposure (e.g., ingestion, inhalation, dermal contact) of a noncarcinogen can determine the type and degree of toxicity. Table 4-3 presents the route-specific RfDs (oral ingestion and inhalation) and the critical toxic endpoints for the chemicals of concern at the EKCO Housewares facility. Dermal RFDs are presented in Table 4-4.

4.3.2.1 Oral Ingestion Route

Chronic oral RfDs were available for most chemicals of concern at the EKCO site. Chemicals for which no oral RfDs were available were benzene, 1,2-dichloroethane, and vinyl chloride. RfDs were not derived for these chemicals.

4.3.2.2 Inhalation Route

Toxicity criteria for inhaled noncarcinogens are now preferentially expressed in IRIS and HEAST as the reference concentration (RfC), rather than the traditional dose of mg/kg-day (RfD). EPA (1993) believes that it is more correct to express toxicity through inhalation based directly on air concentration than by determining internal dose based on pharmacokinetics, surface area adjustments, and the local respiratory effects of sensitizers and irritants. The RfC can usually be converted to an RfD by multiplying by 20m³/day and dividing by 70 kg. Since exposure doses are calculated as mg/kg-day, it is necessary to make this conversion for calculating the hazard quotient.

Oral RfDs were used as inhalation RfDs for volatile organic compounds if the latter were not available from IRIS (1994), HEAST (EPA, 1994) or ECAO. This assumes that an organic chemical producing noncarcinogenic effects by the oral route is likely to produce the same effect through systemic absorption following inhalation (while showering) and that the extent of systemic absorption is comparable through both exposure routes.

Table 4-3 (Cont'd)

Oral and Inhalation Chronic Reference Doses (RfDs) EKCO Housewares Massillon, Ohio

Chemical	Oral RfD (mg/kg-d)	Species/ Critical Effect	Oral RfD Basis/Source ^k	Uncertainty Factor	Inhalation RfD (mg/kg-day)	Species/ Critical Effect	Inhalation RfD Basis/Source ^k	Uncertainty Factor
4-Methyl-2-pentanone	8E-02	rat/kidney and liver effects	13-week gavage study/HEAST	3,000	2E-02	rat/kidney and liver effects	90-day inhalation study/HEAST	1,000
Tetrachloroethene	1E-02	mouse/hepatotoxicity	6-week gavage study/IRIS	1,000	1E-02		Oral RFD	
Toluene	2E-01	rat/liver and kidney effects	13-week gavage study/IRIS	1,000	1.1E-01*	human/neurological effects	occupational exposure/IRIS	300
1,1,1-Trichloroethane	ND				1E+00	guinea pig/liver and lung effects	inhalation study/ECAO	NA
1,1,2-Trichloroethane	4E-03	mouse/clinical serum chemistry	subchronic drinking water study/IRIS	1,000	4E-03		Oral RFD	
Trichloroethene	6E-03	mouse/liver effects	drinking water study/ECAO	3,000	1E-01	human/neurological effects	occupational exposure/ECAO	100
Vinyl chloride	ND				ND			
Xylenes (total)	2E+00	rat/hyperactivity, decreased body weight, increased mortality	chronic gavage study/IRIS	100	2E+100		Oral RFD/	

^{*} Converted from reference concentration (RfC) through the following relationship: RfD (mg/kg-day) = RfC (mg/m³) x 20 m³/day x 1/(70 kg) (20 m³/day is a standard daily inhalation rate from RAGS Supplemental Guidance [U.S. EPA, 1991]).

IRIS - Integrated Risk Information System

HEAST - Health Effects Assessment Summary Tables

ECAO - Environmental Criteria and Assessment Office

ND - Not determined NA - Not available

Table 4-3

Oral and Inhalation Chronic Reference Doses (Rfds) EKCO Housewares Massillon, Ohio

Chemical	Oral RfD (mg/kg-d)	Species/ Critical Effect	Oral RfD Basis/Source ^k	Uncertainty Factor	Inhalation RfD (mg/kg-day)	Species/ Critical Effect	Inhalation RfD Basis/Source ^k	Uncertainty Factor
Acetone	1.0E-01	rat/increased liver and kidney weight and nephrotoxicity	gavage/IRIS	1,000 (100)	1.0E-01		Oral RFD/	****
Benzene	ND				ND			
2-Butanone	6E-01	rat/decreased fetal birth weight	multi-generation developmental study (drinking water)/IRIS	3,000	2.9E-01*	mouse/decreased fetal birth weight	developmental inhalation study/IRIS	3,000
Carbon disulfide	1E-01	rabbit/fetal toxicity, malformations	inhalation teratogenic study/IRIS	100	2.9E-03*	rat/fetal toxicity	inhalation study/HEAST	1,000
Chloroethane	ND				2.9E+00*	mouse/delayed fetal ossification	developmental inhalation study/IRIS	300
Chloroform	1.0E-02	dog/liver lesions	7.5 year oral study/IRIS	1,000	1.0E-02		Oral RFD/	
1,1-Dichloroethane	1E-01	rat/no adverse effects	inhalation study/HEAST	1,000	1.0E-01	cat/kidney damage	13-week inhalation study/HEAST	1,000
1,2-Dichloroethane	ND				ND			
1,1-Dichloroethene	9E-03	rat/hepatic lesions	chronic oral bioassay/IRIS	1,000	9E-03		Oral RFD/	
1,2-Dichloroethene (total)	9E-03	rat/liver lesions	2-year drinking water study/HEAST	1,000	9E-03		Oral RFD/	
Ethylbenzene	1E-01	rat/liver and kidney toxicity	subchronic to chronic oral bioassay/IRIS	1,000	2.9E-01*	rat, rabbit/ developmental toxicity	developmental inhalation studies/IRIS	300

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Table 4-4

Dermal Cancer Slope Factors and Reference Doses
EKCO Housewares
Massillon, Ohio

Chemical	Oral CSF (mg/kg-day) ⁻¹	Dermal CSF ^a (mg/kg-day) ⁻¹	Oral RfD (mg/kg-day)	Dermal RfD ^b (mg/kg/day)
Acetone	NC	NC	1.0E-01	8E-02
Benzene	2.9E-02	3.6E-02	ND	ND
2-Butanone	NC	NC	6E-01	4.8E-01
Carbon disulfide	NC	NC	1E-01	8E-02
Chloroethane	NC	NC	ND	ND
Chloroform	6.1E-03	7.6E-03	1.0E-02	8E-03
1,1-Dichloroethane	ND	ND	1E-01	8E-02
1,2-Dichloroethane	9.1E-02	1.1E-01	ND	ND
1,1-Dichloroethene	6.0E-01	7.5E-01	9E-03	7.2E-03
1,2-Dichloroethene	NC	NC	9E-03	7.2E-03
Ethylbenzene	NC	NC	1E-01	8E-02
4-Methyl-2-pentanone	NC	NC	8E-02	6.4E-02
Tetrachloroethene	5.2E-02	6.5E-02	1E-02	8E-03
Toluene	NC	NC	2E-01	1.6E-01
1,1,1-Trichloroethane	NC	NC	ND	ND
1,1,2-Trichloroethane	5.7E-02	7.1E-02	4E-03	3.2E-03
Trichloroethene	1.1E-02	1.4E-02	6E-03	4.8E-03
Vinyl chloride	1.9E+00	2E+00	ND	ND
Xylenes (total)	NC	NC	2E+00	1.6E+00

NC - Not a carcinogen.

ND - Not determined.

^{*} Oral CSF divided by 0.8 to obtain dermal CSF (80% GI absorption factor).

^b Oral RfD multiplied by 0.8 to obtain dermal RfD (80% GI absorption factor).

4.3.2.3 Dermal Absorption Route

Chronic RfDs have been developed for the oral and inhalation routes, but not for the dermal route. The same criteria and GI absorption factors used for derivation of dermal slope factors were used to develop dermal RfDs for noncarcinogens. A chronic dermal RfD representing the absorbed amount of noncarcinogen was derived by multiplying the value used as the chronic oral RfD by the appropriate GI absorption factor. The GI absorption factor used was 0.8 (volatile organics).

4.3.3 Other Issues Pertaining to Noncarcinogenic Effects

Only chronic RfDs, which are developed to evaluate potential toxicity at greater than seven years of exposure, are presented in Table 4-3 and were used in estimating both childhood and adult noncarcinogenic risk. Subchronic RfDs are sometimes used to evaluate exposures from 2 weeks to 7 years duration. However, chronic RfDs, which are sometimes lower than subchronic RfDs, are used in this risk assessment to ensure a conservative estimate of risk.

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SECTION 5 RISK CHARACTERIZATION

The purpose of the risk characterization is to integrate the information from the exposure assessment and the toxicity assessment in order to estimate quantitatively the potential human health risks that are associated with exposure to the chemicals of potential concern at the EKCO facility. Human health risks are discussed separately for carcinogenic and noncarcinogenic health effects because of the different toxicologic endpoints involved and the different methods that are employed in characterizing the risks. The uncertainties associated with these risk estimates are discussed in Section 6.

Human health risks associated with exposure to carcinogenic chemicals are calculated by multiplying lifetime chemical exposure doses (i.e., daily intakes) by the corresponding chemical-specific cancer slope factors (i.e., risk = daily intake x slope factor). The cancer risk for a particular exposure pathway/route is then estimated by summing the risk estimates for all the chemicals of potential concern through that exposure pathway/route. This approach is in accordance with EPA guidelines on chemical mixtures in which risks associated with carcinogens are considered additive (EPA, 1986). It assumes independence of action by the chemicals (i.e., that there are no synergistic or antagonistic interactions), and that all of the chemicals have the same toxicological endpoint (i.e., cancer). The total cancer risk to an individual member of a receptor population is estimated by summing the combined cancer risks from all relevant exposure routes. The resulting cancer risks are expressed using one significant figure only.

In assessing the carcinogenic risks posed by a site, the NCP establishes a lifetime excess cancer risk of 1E-06 (one-in-one million) as a point of departure for establishing remediation goals. Excess cancer risks lower than 1E-06 are not addressed by the NCP. Excess cancer risks in the range from 1E-06 to 1E-04 (one-in-ten thousand) may or may not be considered acceptable, depending on site-specific factors such as the potential for exposure, technical limitations to remediation alternatives, and data uncertainties. In an

April 1991 OSWER Directive (EPA, 1991) from EPA's Assistant Administrator to Regional Directors, the EPA further clarified the acceptable carcinogenic risk range by stating that when reasonable maximum exposures for both current and future land uses are less than 1E-04 (one-in-ten thousand), action is generally not warranted, unless there are adverse environmental impacts. However, it should be noted that the same directive indicates that the risk manager may decide that a risk less than 1E-04 is unacceptable due to site-specific issues.

Noncarcinogenic health risks are expressed as hazard quotients and hazard indices. The hazard quotient (HQ) for a given chemical in a child or adult is equal to the estimated daily exposure dose (i.e., the daily intake) divided by the acceptable (safe) exposure dose (i.e., the RfD). This approach assumes that exposure to multiple chemicals through a particular exposure route could result in an additive adverse effect, the magnitude of which is proportional to the sum of the HQs. This sum is termed the hazard index (HI). A total HI can also be calculated for an individual receptor that is the sum of the HIs for all relevant exposure routes. When the calculated HQ for an individual chemical or the HI for multiple chemicals exceeds a value of one, there may be a potential for occurence of noncancer health effects. Since the assumption of additivity is most applicable to chemicals that induce the same type of adverse health effect, if the calculated HI is greater than one, then separate HIs can be determined for chemicals according to type of critical effect.

The methodology used to evaluate noncarcinogenic health effects, unlike that used to evaluate carcinogenic effects, does not produce quantitative estimates of risk. If an HQ or HI exceeds one, it simply indicates that there might be a potential for noncarcinogenic health effects to occur under the defined exposure conditions. Conversely, an HQ or HI of less than one indicates that it is unlikely for even sensitive populations to experience adverse noncarcinogenic health effects.

5.1 <u>QUANTITATIVE RISKS ASSOCIATED WITH EXPOSURE TO GROUNDWATER</u> FROM THE UPPER UNIT

Hypothetical future on-site residents using water wells screened in the upper (shallow/intermediate) groundwater unit were assumed to be exposed to site chemicals through drinking water ingestion, noningestion groundwater uses (showering, cooking), and ingestion of garden produce irrigated with groundwater.

5.1.1 Carcinogenic Health Risks

The potential carcinogenic risks associated with hypothetical future residential exposure to groundwater from the upper unit are summarized in Table 5-1. The estimated daily chemical intakes and chemical-specific carcinogenic risks for ingestion uses, noningestion household uses, dermal absorption while bathing, and garden produce ingestion are presented in Appendix B (Tables B-1, B-2, B-3, and B-4, respectively). Total (individual) carcinogenic risk for all exposure pathways was 4E-02 (four-in-one hundred). Inhalation through household noningestion uses contributed about 75% (3E-04) of total risk. The next highest risk was through drinking water ingestion, accounting for 15% (6E-03) of total risk. For the child, most of the risk was through inhalation. The pattern of distribution for the adult was similar to total individual risk. Garden produce ingestion showed a risk of 7E-04, which represented less than 2% of the total risk. Dermal absorption while showering accounted for less than 1% of total risk. Distribution of garden produce ingestion and dermal risk among children and adults showed the source pattern as described for total individual risk.

A summary of those chemicals which exceeded a total individual cancer risk of 1E-06 for the upper unit by exposure pathway is shown in Table 5-2. The two primary compounds contributing to total risk were 1,1-dichloroethene (DCE) (2E-02) and trichloroethane (TCE) (1E-02) through the inhalation pathway. TCE and DCE contributed 4E-03 and 2E-03, respectively, through ingestion of drinking water. Vinyl chloride contributed a total risk of 1E-03 most of which was approximately evenly divided between water ingestion and inhalation. Risks from ingestion of garden produce represented a small percentage of the

Table 5-1

Summary of Potential Carcinogenic Health Risks For the Upper and Lower Units by Exposure Pathway and Age Group Future On-Site Resident

EKCO Housewares, Massillon, Ohio

		Upper Unit				
Exposure Pathway	Individual Cancer Risk	Adult Cancer Risk	Individual Cancer Risk	Child Cancer Risk	Adult Cancer Risk	Individual Cancer Risk
Ingestion of Groundwater	2E-03	4E-03	6E-03	2E-04	4E-04	6E-04
Dermal Absorption while Bathing	1E-04	2E-04	3E-04	7E-06	7E-06	2E-05
Noningestion Groundwater Uses	2E-02	1E-02	3E-02	2E-03	2E-03	3E-03
Ingestion of Garden Produce Irrigated with Groundwater	3E-04	4E-04	7e-04	1E-05	1E-05	3E-05
TOTAL	2E-02	2E-02	4E-02	2E-03	2E-03	4E-03

Table 5-2

Summary of Individual Chemical Cancer Risks Exceeding 1E-06 by Exposure Pathway for the Future On-Site Resident Upper Unit

Upper Unit Exposure Pathway	Chemical	Lifetime Carcinogenic Risk	Percent of Total Lifetime Risk ^e
Drinking Water Ingestion*	Trichloroethene 1,1-Dichloroethene Vinyl chloride 1,2-Dichloroethane 1,1,2-Trichlorethane Tetrachloroethene	4E-03 2E-03 8E-04 1E-05 1E-05 6E-06	9 5 2 0.03 0.03 0.02
Dermal Absorption While Bathing ^b	1,1-Dichloroethane Trichloroethane Vinyl chloride	1E-04 2E-04 2E-05	0.3 0.5 0.04
Noningestion (Inhalation) Uses ^c	1,1-Dichloroethene Trichloroethene Vinyl chloride 1,2-Dichloroethane 1,1,2-Trichlorethane Benzene Chloroform	2E-02 1E-02 6E-04 6E-05 7E-05 3E-06 1E-05	53 26 2 0.2 0.2 <0.01 0.03
Garden Produce Ingestion ^d	Trichloroethene 1,1-Dichloroethene Vinyl chloride	6E-04 1E-04 2E-05	1 0.3 0.05

^{*}Refer to Table B-1 for detailed presentation; Total cancer risk for drinking water ingestion = 6E-03

^bRefer to Table B-2 for detailed presentation; Total cancer risk for dermal absorption = 3E-04

^eRefer to Table B-3 for detailed presentation; Total cancer risk for noningestion inhalation = 3E-02

^dRefer to Table B-4 for detailed presentation; Total cancer risk for garden produce ingestion = 7E-04

^{*}Total lifetime cancer risk for all pathways = 4E-02. Percentages are calculated from risks to two significant figures (Refer to Tables B-1 through B-4).

total, but were associated with risks of 1E-04 for 1,1-DCE and 6E-04 for TCE. Dermal risks were between 1E-05 and 1E-04 for DCE, TCE and vinyl chloride.

The individual cancer risk associated with exposure to 1,1-DCE exceeded 1E-02 for the noningestion (inhalation) exposure pathway and 1E-03 for drinking water ingestion. TCE risk exceeded 1E-02 for inhalation. Garden produce ingestion and dermal absorption while showering accounted for the smallest portion of risk for all chemicals.

The cancer risks associated with groundwater exposure for the upper unit exceeded the range of excess cancer risk (i.e., 1E-06 to 1E-04) set by the NCP for establishing remediation goals.

5.1.2 Noncarcinogenic Health Risks

The potential noncarcinogenic health risks associated with future (child and adult) residential exposure to groundwater in the upper unit are summarized in Table 5-3. The estimated daily chemical intakes and chemical-specific noncarcinogenic risks are presented in Appendix B for drinking water ingestion (Table B-1), dermal absorption while bathing (Table B-2), noningestion (inhalation) uses (Table B-3), and garden produce ingestion (Table B-4).

Table 5-3 shows that total HI for the child was 420, and for the adult was 150. The majority of these potential health effects were through ingestion of drinking water. For the child, HIs were: 240 (drinking water ingestion), 130 (inhalation), 44 (garden produce ingestion), and 11 (dermal absorption during bathing). For the adult, HIs were 100 (drinking water ingestion), 28 (inhalation), 14 (garden produce ingestion), and 5.9 (dermal absorption during bathing).

Table 5-4 presents noncancer risks (HQs) by chemical and exposure pathway for the adult and child. TCE was the primary contributor in all exposure pathways for both the adult and child. The hazard quotients for the child were about 2 to 3 times higher in general than for

Table 5-3

Summary of Potential Noncarcinogenic Health Risks for the Upper and Lower Units by Exposure Pathway and Age Group Future On-Site Resident EKCO Housewares Massillon, Ohio

	Upper Unit		Lower Unit	
Exposure Pathway	Child Hazard Index	Adult Hazard Index	Child Hazard Index	Adult Hazard Index
Ingestion of Groundwater	2.4E+02	1.0E+02	3.8E+00	1.6E+00
Dermal Absorption while Bathing	1.1E+01	5.9E+00	1.5E-01	8.3E-02
Noningestion Groundwater Uses (Inhalation)	1.3E+02	2.8E+01	1.1E+01	2.4E+00
Ingestion of Garden Produce Irrigated with Groundwater	4.4E+01	1.4E+01	5.5E-01	1.7E-01
TOTAL	4.2E+02	1.5E+02	1.6E+01	4.2E+00

the adult. The highest HQ was 240 for TCE for the child via drinking water ingestion. The adult showed a HQ of 100 for TCE. The HQ for the child for dermal absorption of TCE during bathing was 10. TCE was associated with HQs of 44 and 110 for ingestion and inhalation, respectively. Dermal absorption during bathing resulted in the lowest hazard quotients for all chemicals.

Other VOCs that had HQs exceeding one were:

- 1,1-DCE via inhalation
 - 12 (child)
 - 2.5 (adult)
- 1.1-DCE via drinking water ingestion
 - 1.6 (child)
- 1,2-dichloroethane via inhalation
 - 5 (child)
 - 1.1 (adult)
- 1,1,1-trichloroethane via inhalation
 - 2.6 (child)
- 1,1-dichloroethane via inhalation
 - 2.2 (child)
- 1,1,2-trichloroethane via inhalation
 - 1.9 (child)

The noncancer health risks associated with exposure to the upper unit groundwater exceeded unity (1.0) for each exposure pathway examined.

5.2 QUANTITATIVE RISKS ASSOCIATED WITH EXPOSURE TO GROUNDWATER FROM THE LOWER UNIT

Hypothetical future on-site residents using water wells screened in the lower groundwater unit (bedrock aquifer) were assumed to be exposed to site chemicals through drinking water ingestion, inhalation from household noningestion uses, dermal absorption during bathing, and garden produce ingestion via irrigation.

5.2.1 Carcinogenic Health Risks

The potential carcinogenic risks associated with future on-site residential exposure to groundwater from the lower unit are summarized in Table 5-1. The estimated daily chemical intakes and chemical-specific carcinogenic risks are presented in Appendix B for ingestion (Table B-5), dermal absorption while bathing (Table B-6) noningestion (inhalation) household uses (Table B-7), and garden produce ingestion (Table B-8).

Table 5-1 shows that total individual lifetime cancer risk was 4E-03 for the lower unit, which was evenly divided (2E-03) for the child and adult. Inhalation was the greatest total risk of all pathways at 3E-03 (about 75% of total risk for all pathways). Again, this risk was evenly divided between child and adult. Total cancer risk for drinking water ingestion was about 12% of all pathways (6E-04). Dermal absorption while bathing and garden produce ingestion were between 1E-06 and 1E-05, again generally split evenly between child and adult.

Table 5-5 presents a summary for all chemicals by pathway for which total individual cancer risks exceeded 1E-06. The primary contributor to total cancer risk was 1,1-DCE with a risk of 3E-03 (~81% of total). Noningestion use (inhalation) risk for 1,1-DCE was 10 times higher than that for drinking water ingestion. The second highest contributor to total individual risk was vinyl chloride (4E-04; ~12% of total). Risk from drinking water ingestion was approximately equal to risk for noningestion (inhalation) uses.

1,1-DCE was associated with risks of 1E-05, each, for dermal exposure and garden produce ingestion from lower unit groundwater. Vinyl chloride and TCE each produced risks for these pathways between 2E-06 and 6E-06. Total lifetime carcinogenic risks for these three chemicals were approximately two orders of magnitude lower than the risk for these three chemicals by ingestion and inhalation while showering combined (1E-05 vs. 1E-03, respectively). Therefore, both dermal and garden produce ingestion risks were each about 0.1% of total risk for these chemicals.

Table 5-4
Summary of Individual Chemical Noncancer Risks Exceeding A Hazard Quotient of One
By Exposure Pathway and Age Group for the Future On-Site Resident

Upper Unit

		Hazard (Quotient*
Upper Unit Exposure Pathway	Chemical	Adult	Child
Drinking Water Ingestion ^a	Trichloroethene 1,1-Dichloroethene	110 (68) ^f 	240 (56) 1.6 (<1)
Dermal Absorption While Bathing ^b	Trichloroethane	5.9(4)	10(2)
Noningestion (Inhalation) ^c	Trichloroethene 1,1-Dichloroethene 1,1-Dichloroethane 1,2-Dichloroethane 1,1,1-Trichloroethane 1,1,2-Trichloroethane	23 (16) 2.5(2) 1.1(<1) 	110 (26) 12 (3) 2.2 (<1) 5.0 (1) 2.6 (<1) 1.9 (<1)
Garden Produce Ingestion ^d	Trichloroethene	13 (9)	44 (10)

^aRefer to Table B-1 for detailed presentation; Total HI = 100 (adult), Total HI = 240 (child) for drinking water ingestion.

^bRefer to Table B-2 for detailed presentation; Total HI = 5.9 (adult), Total HI = 11 (child) for dermal absorption

^{*}Refer to Table B-3 for detailed presentation; Total HI = 28 (adult), Total HI = 130 (child) for noningestion uses.

dRefer to Table B-4 for detailed presentation; Total HI = 14 (adult), Total HI = 44 (child) for garden produce ingestion.

^{*}Total hazard index for all pathways: Adult = 150; Child = 420.

^{&#}x27;Number in parentheses (next to hazard quotient) is percent of total hazard index for all pathways for the adult or child.

Table 5-5

Summary of Individual Chemical Cancer Risks Exceeding 1E-06

By Exposure Pathway for the Future On-Site Resident

Lower Unit

Lower Unit Exposure Pathway	Chemical	Lifetime Carcinogenic Risk	Percent of Total Lifetime Risk ^e
Drinking Water Ingestion*	1,1-Dichloroethene	3E-04	7
	Vinyl chloride	2E-04	6
	Trichloroethene	4E-05	1
Dermal Absorption While Bathing ^b	1,1-Dichloroethene	1E-05	0.4
	Vinyl chloride	5E-06	0.1
	Trichloroethene	2E-06	0.05
Noningestion (Inhalation) ^c	1,1-Dichloroethene Vinyl chloride Trichloroethene	3E-03 2E-04 1E-04	74 5 3
Garden Produce Ingestion ^d	1,1-Dichloroethene	1E-05	0.4
	Vinyl Chloride	6E-06	0.2
	Trichloroethene	6E-06	0.2

^{*}Refer to Table B-5 for detailed presentation; Total lifetime cancer risk for drinking water ingestion = 6E-04.

^bRefer to Table B-6 for detailed presentation; Total lifetime cancer risk for dermal absorption = 2E-05.

^{&#}x27;Refer to Table B-7 for detailed presentation; Total lifetime cancer risk for noningestion (inhalation) uses = 3E-03.

^dRefer to Table B-8 for detailed presentation; Total lifetime cancer risk for garden produce ingestion = 3E-05.

^{*}Total lifetime cancer risk for all pathways = 4E-03. Percentages are calculated from risks to two significant figures (Refer to Tables B-5 through B-8).

The total cancer risks and certain chemical cancer risks (e.g., 1,1-DCE, TCE and vinyl chloride) associated with the groundwater pathway for the lower unit exceeded the range of excess cancer risk (i.e., 1E-06 to 1E-04) set by the NCP for establishing remediation goals.

5.2.2 Noncarcinogenic Health Risks

The potential noncarcinogenic health risks associated with exposure of a hypothetical future on-site resident to groundwater in the lower unit are summarized in Table 5-3. The estimated daily chemical intakes and chemical-specific noncarcinogenic risks are presented in Appendix B for drinking water ingestion (Table B-5) and noningestion uses (Table B-7). The total hazard indices for the ingestion and noningestion exposure pathways ranged from 4 to 16 for the adult and the child, respectively. The highest HIs were observed for the inhalation pathway (adult, 2.4; child, 11). The second highest contributor to risk by pathway was drinking water ingestion (1.6, adult; 3.8, child). Noncancer HIs from dermal contact and garden produce ingestion represented 30% or less of total HI for the child and adult. Neither of these pathways exceeded an HI of 1 for the adult or child.

The distribution of noncancer effects in the lower unit by chemical is shown in Table 5-6. Three chemicals of all evaluated showed HQs of greater than 1. These were:

- TCE via inhalation
 - 1.1 (child)
- TCE via drinking water ingestion
 - 2.5 (child)
 - 1.1 (adult)
- 1,1-DCE via inhalation
 - 1.7 (child)
- 1,2-DCE via inhalation
 - 7.6 (child)
 - 1.6 (adult)

Table 5-6
Summary of Individual Chemical Noncancer Risks Exceeding A Hazard Quotient of One
By Exposure Pathway and Age Group for the Future On-Site Resident

Lower Unit

Upper Unit Exposure Pathway	Chemical	Hazard Adult	Quotient" Child
Drinking Water Ingestion	Trichloroethene	1.1 (26) ^d	2.5 (16)
Noningestion (Inhalation) ^b	Trichloroethene 1,1-Dichloroethene 1,2-Dichloroethene	 1.6 (38)	1.1 (7) 1.7 (11) 7.6 (48)

^aRefer to Table B-5 for detailed presentation; Total HI = 1.6 (adult), Total HI = 3.8 (child) for drinking water ingestion.

^bRefer to Table B-7 for detailed presentation; Total HI = 2.4 (adult), Total HI = 11 (child) for noningestion uses.

[&]quot;Total hazard index for all pathways: Adult = 4.2; Child = 16.

dNumber in parentheses (next to hazard quotient) is percent of total hazard index for all pathways for the adult or child.

The noncancer health risks for the child and for the adult associated with the lower groundwater unit exceeded unity (1.0) for the drinking water ingestion exposure pathway and for the noningestion (inhalation) exposure pathway.

5.3 CONCLUSIONS

Total cancer risks in the upper unit were 4E-02. The major contributors (i.e., greater than 1E-04) to risk were from TCE, 1,1-DCE and vinyl chloride through drinking water ingestion and inhalation while showering and for TCE for garden produce ingestion. Total dermal cancer risks were 3E-04 with TCE exceeding a risk of 1E-04. Total HI by all pathways for the child was 420, and for the adult was 150. Total HI exceeded 1 for each exposure pathway for the child and adult, with drinking water ingestion producing the majority of risk. Garden produce ingestion was about 10% of total HI (adult and child) and dermal HI from 1 to 3%. TCE and 1,1-DCE were the principal contributors to hazard index primarily through drinking water ingestion and inhalation. TCE contributed approximately 5 to 10% of total HI for garden produce ingestion (adult and child, respectively) and less than 5% for dermal absorption while bathing.

For the lower unit, the total cancer risk (4E-03) was an order of magnitude lower than that from the upper unit. About 75% of total lifetime cancer risk was associated with inhalation from household uses (bathing, cooking), and was split evenly between childhood (ages 1 to 6 years) and adult (ages 7 to 30 years) exposure. The majority of this risk was due to the inhalation of 1,1-DCE, TCE and vinyl chloride. About 15% of total risk was due to drinking water ingestion, and about 1% each were due to garden produce ingestion and dermal absorption while bathing. The drinking water ingestion risk was influenced primarily by 1,1-DCE and vinyl chloride.

For noncancer risks in the lower unit, the majority of risk was associated with inhalation from household noningestion uses for both the child (approximately 70%) and adult (approximately 60%). 1,2-DCE contributed to the greatest extent through this exposure pathway for both the child and adult. Drinking water ingestion (primarily TCE) also

contributed substantially to child and adult risk, with HQs greater than 1.

Future exposure to the lower groundwater unit posed less total cancer and noncancer risks than to the upper unit. Total risks for both units were greater than the points of departure of 1E-04 (cancer risk) and a hazard index of one (noncancer risk). A discussion of uncertainties and assumptions is presented in the next Section (6) of the report.

5.4 <u>REFERENCES</u>

EPA (United States Environmental Protection Agency), 1986. Guidelines for Carcinogenic Risk Assessment. Federal Register. 51 (185):33992.

EPA (United States Environmental Protection Agency), 1991. Role of the Baseline Risk Assessment in Superfund Remedy Selection Decisions. Memorandum from Don R. Clay, Assistant Administrator. OSWER Directive 9355.0-30.

SECTION 6 UNCERTAINTY ANALYSIS

6.1 <u>INTRODUCTION</u>

The principal goals of the uncertainty analysis are to provide to the appropriate decision makers a discussion of the key assumptions made in the risk assessment that significantly influence the risk results and to assess the contribution of these factors to the under- or overestimation of risk. The uncertainty analysis should show that the calculated risks are relative in nature and do not represent an absolute quantification.

In recent months, the U.S. EPA has placed even more emphasis on the uncertainty analysis. In a 26 February 1992 memorandum from the Deputy Administrator to all assistant and regional administrators (EPA, 1992), U.S. EPA provides additional guidance on explaining risks and all their underlying data so that the strengths and weaknesses of the assessment become clear. This section of the risk assessment attempts to explain the key assumptions used in this report and to present the range of variability inherent in these assumptions.

In the absence of empirical or site specific data, assumptions are developed based on best estimates of data quality, exposure parameters and dose-response relationships. To assist in the development of these estimates, the EPA recommends the use of guidelines and standard exposure factors in risk assessments conducted under CERCLA (EPA, 1989; 1991). The use of these standard factors is intended to promote consistency among risk assessments where assumptions must be made for high-end (RME) risk estimates. Although the use of standard factors no doubt promotes comparability, their usefulness in accurately predicting risk is directly proportional to their applicability to the site-specific conditions.

The carcinogenic and noncarcinogenic risk estimates for the EKCO Facility were based on a number of assumptions that incorporated varying degrees of uncertainty resulting from several sources, including those pertaining to:

- Selection of exposure pathways, input parameters, algorithms and scenarios.
- Confidence in toxicological data used to estimate cancer potency factors and noncancer reference doses.

6.2 <u>SUMMARY OF SIGNIFICANT CANCER AND NONCANCER RISKS AT THE</u> EKCO HOUSEWARES FACILITY

Tables 6-1, 6-2 and 6-3 summarize the chemicals and exposure pathways that contributed greater than 1E-04 to carcinogenic risk and greater than a hazard quotient of 10 to noncarcinogenic risk in the upper and lower groundwater units for the hypothetical future on-site resident.

6.2.1 Upper Unit Groundwater Risks

Table 6-1 shows that of all chemicals evaluated, the primary contributor to total cancer risk in the upper unit (4E-02) was 1,1-DCE with a chemical risk of 2E-02 (~60% of total) through inhalation from household noningestion uses. TCE contributed about 35% (approximately 1E-03) of total risk, the major portion of this risk being through inhalation noningestion household uses. Vinyl chloride presented about 3% (1E-04) of total risk chiefly through drinking water ingestion and inhalation from household noningestion uses.

Table 6-2 shows that TCE was the primary chemical producing potential noncancer health effects, contributing about 92% of the total child HI, and about 93% of the adult HI. This amounted to an HQ of about 400 and 140 for the child and adult, respectively. About 23% of total child HI (HQ=12) was accounted for by 1,1-DCE.

6.2.2 Lower Unit Groundwater Risks

Table 6-3 shows that two chemicals in the lower unit groundwater exceeded cancer risks of 1E-04: 1,1-DCE and vinyl chloride. The highest contributor was 1,1-DCE via inhalation through household noningestion uses (about 74%). No chemicals exceeded a noncarcinogenic hazard quotient of 10 for any exposure pathway evaluated in the lower unit.

Table 6-1
Summary of Individual Chemical Cancer Risks Exceeding 1E-04 by Exposure Pathway for the Future On-Site Resident

Upper Unit

Upper Unit Exposure Pathway	Chemical	Lifetime Carcinogenic Risk	Percent of Total Risk ^e
Drinking Water Ingestion*	Trichloroethene 1,1-Dichloroethene Vinyl chloride	4E-03 2E-03 8E-04	9 5 2
Dermal Absorption While Bathing ^b	Trichloroethene	2E-04	0.5
Noningestion (Inhalation) ^c	1,1-Dichloroethene Trichloroethene Vinyl chloride	2E-02 1E-02 6E-04	53 26 2
Garden Produce Ingestion ^d	Trichloroethene	6E-04	1

^aRefer to Table B-1 for detailed presentation; Total cancer risk for drinking water ingestion = 6E-03

^bRefer to Table B-2 for detailed presentation; Total cancer risk for dermal absorption = 3E-04

^{&#}x27;Refer to Table B-3 for detailed presentation; Total cancer risk for noningestion (inhalation) = 3E-02

dRefer to Table B-4 for detailed presentation; Total cancer risk for garden produce ingestion = 7E-04

^{*}Total lifetime cancer risk for all pathways = 4E-02

Table 6-2
Summary of Individual Chemical Noncancer Risks Exceeding A Hazard Quotient of Ten
By Exposure Pathway and Age Group for the Future On-Site Resident

Upper Unit

Upper Unit Exposure Pathway	Chemical	Hazard (Adult	Quotient ^d Child
Drinking Water Ingestion ^a	Trichloroethene	100 (68)°	240 (56)
Noningestion (Inhalation) ^b	Trichloroethene 1,1-Dichloroethene	23 (16)	110 (26) 12 (3)
Garden Produce Ingestion ^c	Trichloroethene	13 (9)	44 (10)

⁸Refer to Table B-1 for detailed presentation; Total HI = 100 (adult), Total HI = 240 (child) for drinking water ingestion.

^bRefer to Table B-3 for detailed presentation; Total HI = 28 (adult), Total HI = 130 (child) for noningestion (inhalation).

Refer to Table B-4 for detailed presentation; Total HI = 14 (adult), Total HI = 44 (child) for garden produce ingestion.

^dTotal hazard index for all pathways: Adult = 150; Child = 420.

[&]quot;Number in parentheses (next to hazard quotient) is percent of total hazard index for all pathways for the adult or child.

Table 6-3
Summary of Individual Chemical Cancer Risks Exceeding 1E-04
By Exposure Pathway for the Future On-Site Resident

Lower Unit

Lower Unit Exposure Pathway	Chemical	Lifetime Carcinogenic Risk	Percent of Total Lifetime Risk ^c
Drinking Water Ingestion ^a	1,1-Dichloroethene	3E-04	7
	Vinyl chloride	2E-04	6
Noningestion (Inhalation) ^b	1,1-Dichloroethene	3E-03	74
	Vinyl chloride	2E-04	5

^aRefer to Table B-5 for detailed presentation; Total lifetime cancer risk for drinking water ingestion = 6E-04.

^bRefer to Table B-7 for detailed presentation; Total lifetime cancer risk for noningestion (inhalation) = 3E-03.

Total lifetime cancer risk for all pathways = 4E-03.

6.2.3 Conclusions

These chemicals and exposure pathways, which were the highest contributors to the site risks, and their associated uncertainties are the main focus of the following discussion. General uncertainties inherent in the risk assessment process (i.e., standard default assumptions) are presented where they are relevant to the EKCO Facility.

6.2.4 Uncertainties Associated with Exposure Assessment

The exposure assumptions directly influence the calculated doses (daily intakes), and ultimately the calculation of risk. In general, conservative exposure assumptions were made in calculating exposure doses such as the selection of exposure routes and scenarios, and the exposure input factors (e.g., contact rate, exposure frequency, exposure duration, body weight and surface area) used to estimate exposure doses. In most cases, these uncertainties contribute to overestimation of plausible real-life exposures, and therefore, true risk is overestimated. These assumptions are appropriate when performing risk assessments of this type so that risk managers can be reasonably assured that human health risks are not underestimated, and so that risk assessments for different sites can be compared.

The concept of reasonable maximum exposure (RME) was used to develop exposure doses for the hypothetical future on-site resident, and is defined as the "maximum exposure that is reasonably expected to occur at the site (EPA, 1989)." Several significant variables that determine the exposure doses for the RME are based on upper-bound (typically 90th to 95th percentile or greater) estimates. These are:

- The 95% upper confidence limit of the average medium concentration for the chemical used to calculate the exposure dose.
- Intake/contact rate (IR) (upper-bound value).
- Exposure frequency (EF) (upper-bound value).
- Exposure duration (ED) (upper-bound value).

Therefore, the calculated exposure dose for any given chemical, which results from integration of all of these variables, represents an upper-bound (high-end) estimate of the probable exposure dose. The use of these upperbound exposure parameters, coupled with conservative estimates of toxicity, in turn will yield risk results that represent an upper-bound estimate of the carcinogenic and noncarcinogenic health risks. Below are discussed several site specific uncertainties which relate to calculation of exposure concentration.

6.2.4.1 Improbable Use of Upper Unit as a Drinking Water Source

Exposure to the upper groundwater unit is highly improbable in the future in view of the fact that this unit is not currently used because of its low yield (See Appendix C). The majority of both cancer and noncancer risks calculated for this unit was through ingestion and household noningestion (inhalation while showering) uses. Relatively small percentages of total risk were associated with garden produce ingestion and dermal contact during bathing for this unit.

In view of the very low probability that the upper unit would ever be used as a drinking water source in the future, the risks calculated for the upper unit based on drinking water ingestion and household noningestion uses should be viewed as hypothetical. Assuming the upper unit was used in the future as a source for watering of home grown vegetables, total lifetime cancer risk for the upper unit would be about 7E-04 and the noncancer HI would be about 44 (child) and 14 (adult).

6.2.4.2 Dilution and Degradation of Volatiles in Upper and Lower Units

Exposure concentrations of VOCs calculated for both the upper and lower groundwater units are likely overestimated to a high degree relative to a 30-year residence time by future on-site residents. It was conservatively assumed in the risk calculations that there would be no degradation or dilution of the VOCs in groundwater.

Also, it is noted that during watering of gardens, VOCs in the groundwater will volatize, and therefore the concentration of VOCs that was used to estimate vegetable and fruit uptake is overestimated. VOC concentrations are expected to decrease significantly during the 30 year lifetime of the future resident. TCE, for example, evaporates from water very quickly (minutes to hours), and although degradation processes are slow, the approximate half-life of TCE in water influenced by biodegradation is about 10 to 11 months (Howard, 1990). Moreover, based on pump and treatment systems currently in operation, removal of VOCs over a period of years will result in relatively lower exposure concentrations averaged over the next 30 years. This expected reduction in the estimated exposure point concentration will be proportionately reflected in lower cancer and noncancer risks associated with both the upper and lower groundwater units.

6.2.4.3 Central Tendency Exposure Issues

The "Habicht Memorandum" (EPA, 1992) on risk characterization recommends a discussion of the range of exposures and multiple risk descriptors in the uncertainty analysis (e.g., central tendency) in addition to the RME calculations, which are consistent with the mandate to evaluate the "high end" risk descriptor (EPA, 1992). Central tendency risk descriptors are described by EPA (1992) as "either the arithmetic mean risk or the median risk." RME is defined as "the highest exposure that is reasonable expected to occur at a site" (EPA, 1989). In practice, the RME is estimated by combining upper bound (90-95th percentile) values for some but not all exposure parameters. WESTON obtained relevant central tendency exposure values from EPA Region V (Personal Communication, 1994). These were used to estimate average risks for those chemicals having significant impact on the (RME) risk presented for the upper and lower groundwater units (Section 5).

Table 6-4 compares the exposure input values used to obtain RME and average risk. The central tendency values were incorporated into the exposure algorithms in Section 3 where appropriate (Table 3-1, drinking water ingestion; Table 3-2, inhalation while bathing), and Table 3-3, ingestion of garden produce). Dermal absorption was not evaluated since this pathway was a minor contributor to total risk. A ratio of the central (CT) tendency

Table 6-4

Comparison of Exposure Input Values^a for

RME and Average (Central Tendency) Exposure Dose Calculations

	Input Values						
Variable	RME		Central Tendency				
	Child	Adult	Child	Adult			
IR (m³/day)	15	15	15	15			
IR (L/day)	1	2	0.7	1.4			
EF (days/year)	350	350	234	234			
ED (years)	6	24	2	7			
AT-Noncancer (days)	2,190	8,760	730	2,555			
AT-Cancer (days)	25,550	25,550	25,550	25,550			

RME = Reasonable maximum exposure

IR = Ingestion rate (drinking water)

EF = Exposure frequency ED = Exposure duration

AT = Averaging time (cancer, or noncancer)

^a Only those exposure input variables shown in the table were changed. Inhalation rates for showering are shown but did not change in the inhalation exposure equation. All other values (Tables 3-1, 3-2, 3-3, and 3-4) remained identical (i.e., the RME default variables).

exposure to the RME exposure was calculated for each exposure algorithm for the child and adult (excluding the "C_w" term). The detailed calculations are presented in Table B-9 in Appendix B. This ratio was then multiplied with the respective RME risks to obtain the average risk. The results are summarized in Table 6-5. The range of both child and adult cancer and noncancer risks decreased by 30 to 90% compared to the RME risk. The decrease was dependent on the groundwater unit, chemical and exposure pathway compared to the RME risk.

6.2.5 <u>Uncertainties Associated with Toxicity Assessment</u>

For a risk to exist, both significant exposure to the pollutants of concern and toxicity at these predicted exposure levels must exist. The toxicological uncertainties primarily relate to the methodology by which carcinogenic and noncarcinogenic criteria (i.e., cancer slope factors and reference doses) are developed. In general, the methodology currently used to develop cancer slope factors and reference doses is conservative, and likely results in overestimation of human toxicity (EPA, 1989). These and other factors are discussed in the subsections below.

6.2.5.1 Cancer Slope Factors

Although there is evidence to suggest some carcinogens may exhibit thresholds, cancer slope factors are developed assuming there is no safe level of exposure to any chemical proven or suspected to cause cancer. This uncertainty implies that exposure to even a single molecule of a chemical may be associated with a finite risk, however small. The assumption is that even if relatively large doses of a pollutant were required to cause cancer in laboratory animals (i.e., much higher than a person would ever likely be exposed to over a lifetime), these exposure doses can be linearly extrapolated downward many orders of magnitude to estimate slope factors for humans. A significant uncertainty for the carcinogens is whether the cancer slope factors accurately reflect the carcinogenic potency of these chemicals at low exposure concentrations. The calculated slope factor is used to estimate an upperbound lifetime probability of an individual developing cancer as a result

Table 6-5

Comparison of Average and RME Risk Estimates^a for Selected Chemicals in the Upper and Lower Groundwater Units

Groundwater Unit	Health Effect Evaluated	Chemical	Exposure Pathway	RME Risk		Average Risk	
				Child	Adult	Child	Adult
Upper	Cancer	TCE	Water Ingestion	1.3E-03	2.3E-03	2.0E-04(85)b	3.1E-04 (87)
		1,1-DCE	Inhalation	1.1E-02	9.3E-03	2.5E-03 (77)	1.8E-03 (81)
	Noncancer	TCE	Water Ingestion	240	100	110 (54)	47 (53)
		TCE	Inhalation	110	23	74 (33)	15 (35)
		TCE	Vegetable Ingestion	44	13	29 (34)	8.7 (33)
Lower	Cancer	1,1-DCE	Water Ingestion	1.0E-04	1.8E-04	1.6E-05 (84)	2.5E-05 (86)
		1,1-DCE	Inhalation	1.5E-03	1.3E-03	3.3E-04 (78)	2.5E-04 (81)

^{*} Refer to Table B-9 in Appendix B for detailed calculations.

^b Numbers in parentheses refer to percentage decrease in converting from RME to average risk.

of exposure to a particular level of a carcinogen. Therefore, the cancer slope factors developed by EPA are generally conservative and represent the upperbound limit of the carcinogenic potency of each chemical. The actual risk posed by each chemical is unknown, but is likely to be lower than the calculated risk, and may even be as low as zero (EPA 1989). The conclusion is that these toxicity assumptions will typically result in an overestimation of carcinogenic risk.

The assumption that all carcinogens (whether A, B1, B2, or C) can cause cancer in humans is also conservative. Only those chemicals classified as "A" carcinogens by the EPA are unequivocally considered human carcinogens. The other three classes are <u>probable</u> (B1, B2) or <u>possible</u> (C) human carcinogens. In this risk assessment, all "probable" and "possible" carcinogens are given the same weight in the toxicity assessment (and consequently in the estimation of risk) as true human carcinogens. This assumption most likely overestimates actual carcinogenic risk to the human.

1,1-Dichloroethene (1,1-DCE)

Although 18 studies have been evaluated by EPA for potential cancer risk, only a single inhalation animal toxicity study is considered by EPA to be sufficient evidence that 1,1-DCE is a complete carcinogen (ATSDR, 1989). Thus, there is question as to whether 1,1-DCE is truly an inhalation carcinogen in humans. It is currently classified as a "C" carcinogen (see Table 4-2). Therefore, the potential risk from exposure by inhalation (noningestion household uses) and ingestion of groundwater may be moderately overestimated.

Trichloroethene

The oral and inhalation slope factors for trichloroethene (TCE) are currently under review by EPA (IRIS, 1994). Until questions regarding the adequacy of the critical studies and the toxicity criteria for TCE are resolved, there is some uncertainty in estimating its potential cancer risk. Since the slope factor is subject to change, it is not possible to calculate accurately the cancer risk associated with TCE.

Vinyl Chloride

Currently, vinyl chloride is under review by the EPA (IRIS, 1994). A value of 1.9E+0 (mg/kg-day)⁻¹, which is based on scientific data from 1984 and 1985, was used for the oral cancer slope factor (EPA, 1994). The slope factor is likely to change soon in view of recently available data, but which has not yet been completely evaluated by U.S. EPA. In their view (IRIS, 1994) the potency of vinyl chloride as an oral carcinogen is likely to increase. Consequently, the cancer risk due to the ingestion of vinyl chloride may be underestimated.

6.2.5.2 Reference Doses

In the development of reference doses (RfDs) for each chemical by exposure route, it is assumed that a threshold dose exists below which there is no potential for adverse health effects to the most sensitive individuals in the population. The RfD is typically derived from dose-response studies in animals in which a NOAEL (no-observed-adverse-effect level) or a LOAEL (lowest-observed-adverse-effect level) is determined by applying several uncertainty factors of 10 each. An additional modifying factor of up to 10 can be applied which accounts for a qualitative professional assessment of additional uncertainties in the available toxicity data (EPA, 1989). The final degree of extrapolation for a given chemical can range anywhere between 10 and 100,000 and therefore result in a human subthreshold dose of one tenth to one-hundred thousandth of the study dose. In general, the calculated RfD is overly protective, and its use probably results in a moderate to high overestimation of noncarcinogenic risk.

Use of Chronic RfDs in Children

Oral chronic RfDs were used in calculating hazard quotients for the 1 to 6 year old child. The use of chronic RfDs in this age group is conservative and will result in an overestimation of risk. Chronic RfDs are developed assuming a lifetime daily exposure. Subchronic RfDs, which are usually based on an exposure duration of 2 weeks to 7 years,

generally tend to be greater than chronic RfDs, and therefore, would in some cases result in a lower hazard quotient and index.

6.2.6 Summary and Conclusions of Uncertainy Analysis

Cancer risks based on future residential development in both upper and lower units exceeded the 1E-04 benchmark used by EPA (1992a) with the upper (1E-02) being about one order of magnitude higher than the lower unit (1E-03). The same pattern was evident for noncancer risks. For both cancer and noncancer risks, the largest contributors by chemical were TCE, 1,1-DCE and vinyl chloride, with groundwater ingestion and inhalation while showering posing the greatest risks.

Table 6-6 summarizes the primary uncertainties and assumptions used to evaluate risks at the EKCO Housewares Facility and their relative impact on risk. It is concluded from this uncertainty analysis that site risk has been moderately to highly overestimated based on an integration of all key assumptions used in the risk assessment for this facility.

Several important factors at the EKCO Housewares Facility support their conclusion:

- Based on the unlikely use of the upper unit for drinking water and household noningestion uses, the total risks for this unit are significantly overestimated.
 The total risks associated with this unit are most relatistically associated with use of this unit for watering of home grown vegetables and fruits.
- There is a low probability that the site will be developed residentially, and therefore risks for both units are likely overestimated.
- Due to dilution and degradation processes and pump and treatment systems already in place, the actual exposure concentration to all VOCs are expected to be significantly lower averaged over the next 30 years.

Table 6-6

Summary of Uncertainty Analysis EKCO Housewares Facility

		Effects on Risk Esti	mate
Uncertainty Element	Potential for Overestimation	Potential for Underestimation	Potential for Over- or Underestimation
Exposure Assessment			
Standard assumptions regarding body weights, skin surface areas, inhalation rates, and life expectancy			Low
Media intake rates	Moderate		
Exposure frequencies	Moderate		
Exposure durations	Moderate		
Use of upper unit for residential drinking water and other household uses	High		
Dilution and degradation of VOCs	High		
Toxicity Assessment	. —		
Use of chronic RfDs for estimating noncancer risk in children	Moderate		
Cancer slope (potency) factors	High		
1,1-DichloroethyleneTrichloroethyleneVinyl chloride	Moderate	Moderate	High
Reference doses	Moderate-High		·

6.3 **REFERENCES**

ATSDR (Agency for Toxic Substances and Disease Registry), 1989. *Toxicological Profile for 1,1-Dichloroethethylene*, Final Report, Clement Associates, Contract No. 205-88-0608, PB90-182114.

EPA (U.S. Environmental Protection Agency), 1989. Risk Assessment Guidance for Superfund, Volume 1, Human Health Evaluation Manual (Part A). Interim Final. Office of Solid Waste and Emergency Response, Washington, D.C. EPA/540/1-89/002.

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Howard, P.H., 1990. Handbook of Environmental Fate and Exposure Data for Organic Chemicals, Volume II, Solvents. "Trichloroethylene." p. 467. Lewis Publishers, Chelsea, MI

IRIS (Integrated Risk Information System), 1994. Computerized data base of toxicity information or hazardous chemicals maintained by U.S. EPA, Washington, D.C.

Personal Communication. 1994. Phone conversation between Sally Averill, Project Manager, EPA Region V, RCRA Branch, and Robert Warwick, Ph.D., Technical Manager, Risk Assessment, Roy F. Weston, Co. 8 November 1994.

APPENDIX A
RAW DATA

Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS. HSL LIST

pri p			VOI.	ATILES BY	GC/M	Coast La S, HSL LI	ST		Report Dat	:e: (05/20/ 94 11:1
RFW Batch Number: 9405G855	<u> </u>	Client: [Ekco	Houseware	<u> </u>		Work	Order: 0299	94-002-005-9		Page: 1
	Cust ID:	L-1		L-1	l	L-1 Dup	•	L-1 Dup	L-a	2	L-2
Sample	RFW#:	001		001 DI	-	002	2	002 DL	003)	003 DL
Information	Matrix:	WATER		WATER		WATER		WATER	WATER		WATER
	D.F.:	1.0)		. 0	1.		5.0	1.	.0	20.0
	Units:	ug/L		ug/l	-	ug/l	-	ug/L	ug/l	-	ug/L
	uene-d8	98	%	102	%	96	%	106 9	6 95	%	
Surrogate Bromofluoro		96 ⁻	%	110	%	96	%	107 9	6 99	%	110 %
Recovery 1,2-Dichloroet		96	%	111	%	97	%	111 9		%	114 %
	****	========			-=f]==				[]====== = ===	f1-	======================================
Chloromethane		2	U	NA		2	U	NA	2	U	NA
Bromomethane		2	U	NA		2	U	NA	2	U	NΛ
Vinyl chloride		20		NA		19		NA	1	U	NΛ
Chloroethane		2	U	NA		2	U	NA	2	U	NA
Methylene Chloride		1	U	NA		1	U	NA	1	U	NΛ
Acetone		7	В	NA		6	В	NA	4	В	NA
Carbon disulfide		1	U	NA		1	U	NA	1	U	NΛ
1,1-Dichloroethene		2		NA		2		NA	1	U	NA
1,1-Dichloroethane			E	26			E	31	1	U	NΛ
1,2-Dichloroethene (total)			E	42			E	42	2		NA
Chloroform		1	U	NA		1	U	NA	ĺ	U	NA
1,2-Dichloroethane		1	U	NA		1	U	NA	1	U	NΛ
2-Butanone		2	U	['] NA		2	U	NA	2	Ũ	NA
1,1,1-Trichloroethane		4		NA		4		NA	5	•	NA
Carbon Tetrachloride		i	U	NA		ì	U	NA	ĭ	U	NA
Vinyl acetate		2	Ŭ	NA		2	Ū	NA	2	Ū	NA
Bromodichloromethane		ĩ	Ü	NA		ī	Ŭ	NA	ī	Ū	NA
1,2-Dichloropropane		i	Ū	NA		ī	Ũ	NA	i	Ũ	NA
air 1 2 Diebleverse		i	Ū	NA		ī	Ū	NA	ī	Ū	NA
Trichloroethene		-	Ě	42		_	Ē	41	_	Ē	240
Dibromochloromethane		1	Ū	NA		1	Ũ	NA	1	Ū	NA
1,1,2-Trichloroethane		ī	Ū	NA		Ī	Ū	NA	i	Ū	NA
Benzene		i	Ŭ	NA		ī	Ŭ	NA	i	Ŭ	NA
trans-1,3-Dichloropropene		i	Ŭ	NA		ī	Ŭ	NA	i	Ū	NA
Bromoform		i	Ŭ	NA		i	Ŭ	NA	i	Ŭ	NA
4-Methy1-2-pentanone		ż	Ŭ	NA		Ž	Ũ	NA	ż	Ŭ	NA
2-Hexanone		ž	Ŭ	NA		ž	Ŭ	NA	2	Ŭ	NΛ
Tetrachloroethene		ĩ	Ŭ	NA		ĩ	Ŭ	NA	ī	Ŭ	NΛ
1,1,2,2-Tetrachloroethane		i	Ŭ	NA		i	Ŭ	NA	i	Ŭ	NΛ
*= Outside of EPA CLP QC	imits.	•	•			-	-		-	-	

RFW Batch Number: 9405G855	Client: Ekco	Houseware	Work	Order: 02994-0	02-005-9	Page: 1b
Cust ID:	L-1	L-1	L-1 Dup	L-1 Dup	L-2	L-2
REW#:	001	001 DL	002	002 DL	003	003 DL
Toluene Chlorobenzene Ethylbenzene Styrene Xylene (total) *= Outside of EPA CLP QC limits.	1 U 1 U 1 U 1 U	NA NA NA NA	1 U 1 U 1 U 1 U	NA NA NA NA NA	1 U 1 U 1 U 1 U	NA NA NA NA NA

±0-0 -

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Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATHES BY GC/MS. HSL LIST

RFW Batch Number: 9405G855	Client: Ek	VOLATILES BY CO Houseware	GC/MS, HSL		k Order: 02	994-	Report Date -002-005-9	e:	05/20/94 11:11 Page: 2a
Cust ID	: L-3	L-4	t	4	L-4		L-5		L-5
Sample RFW# Information Matrix D.F.	: WATER : 1.0	005 WATER 1.	WATE 0	R 1.0	005 MSD WATER 1.	0	006 WATER 1.0	0	006 DL WATER 10.0
Units	: ug/L	ug/L	uç	g/L	ug/L		ug/L		ug/L
Toluene-de Surrogate Bromofluorobenzen Recovery 1,2-Dichloroethane-de	e 99 9 4 99 9	6 99 6 95	% 93 % 103 % 93	1 %	96 103 99	% %	103 106 111	% %	97 % 101 % 98 %
Chloromethane	2 (J 2	Ü	2 U	2	≖f]. U	2	-f1 U	NA I
Bromomethane Vinyl chloride		· -	U U	2 U 1 U	2	U	2	ñ	NA E 1
Chloroethane	- 1 C	· · · · · · · · · · · · · · · · · · ·	U	2 U	2	U	2	E	51 NA
Methylene Chloride	— เ้เ	•	Ŭ	īŬ	ī	Ŭ	ī	Ŭ	NA
Acetone		3	В	3 B	Ž	Ŭ	16	B	NA
Carbon disulfide		1	U	1 U	1	U	1	U	NA
1,1-Dichloroethene		J 1	U 10	l %	100	%	1	U	NA
1,1-Dichloroethane	1 (,		5	5		18		NA
1,2-Dichloroethene (total)	<u> </u>			9	20		33		NA
Chloroform		•	U	1 U	1	U	1	U	NA
1,2-Dichloroethane		J 1	U	1 U	1	Ü	1	U	NA
2-Butanone	_ 2 (] 2	U	2 U	2	Ü	2	Ü	NA
1,1,1-TrichToroethane	_ !!	•	U	1 U	I .	U	1	U	NA
Carbon Tetrachloride	_ 1 !	-	์ บ - บ	1 U	1	U	1	Ü	NA
Vinyl acetate Bromodichloromethane		•	U	1 11	2	U	2	U	NA NA
1,2-Dichloropropane	i i	•	Ü	1 0	1	U	1	U	NA NA
cis-1,3-Dichloropropene	– i i	•	Ŭ	i Ü	i	Ü	1	Ü	NA NA
Trichloroethene	– i i	•	Ü 9		98	%	1	Ŭ	NA NA
Dibromochloromethane	– i i	•	Ü	່າ ບິ	701	ũ	i	Ŭ	NA NA
1,1,2-Trichloroethane	– ii	•	Ŭ	iŭ	i	Ŭ	i	Ŭ	NA
Benzene	– i i	-	Ü 104	_	106	% .	i	Ū	NA
trans-1,3-Dichloropropene	– ii	j	Ü	1 Ü	1	Ũ	ī	Ū	NΛ
Bromoform	i i	J	U	1 U	1	U	1	U	NA
4-Methy1-2-pentanone	2 l		U	2 U	2	U	2	U	NA
2-Hexanone	_ 2 (•	U	2 U	2	U	2	U	NA
Tetrachloroethene	1 (U	1 U	1	U	1	U	ĦΛ
1,1,2,2-Tetrachloroethane *= Outside of EPA CLP QC Timits.	1 (J 1	U	1 U	1	U	1	U	ил

RFW Batch Number: 9405G855	Client: Ekco H	ouseware	Work	Order: 02994-00	2-005-9	Page: 2b
Cust ID:	L-3	L-4	L- 4	L-4	L-5	L-5
RFW#:	004	005	005 MS	005 MSD	006	006 DL
Toluene	1 U		94 %	95 %	1 U	NA
Chlorobenzene	1 U	1 U	96 %	97 %	1 U	NΛ
Ethylbenzene	1 U	1 U	, 1 U	1 U	1 U	NA
Styrene	1 U	l Ú	, 1 U	וט	1 0	NA
Xylene (total) *= Outside of EPA CLP QC limits.	1 U	1 U	1 U	1 V	in 1 U	NA

AHP On-Site Shallow Wells Minimum and Average DTWs - 1994 10/28/94

				2/15/94		5/3/94		8/10/	DTW DOL	
Well	TOC	GL	Stick- Up	DTW TOC	DTW BGL	DTW TOC	DTW BGL	DTW TOC	DTW BGL	DTW BGL Arithmetic Mean
L-1	946.33	944.2	2.13	20.68	18.55	20.75	18.62	24.65	22.52	19.90
L-2	947.57	946.2	1.37	17.49	16.12	15.38	14.01	17.11	15.74	15.29
L-3	946.91	946.0	0.91	15.25	14.34	14.54	13.63	16.52	15.61	14.53
L-4	938.22	935.9	2.32	8.28	5.96	8.08	5.76	8.45	6.13	5.95
L-5	936.98	934.7	2.28	7.67	5.39	7.39	5.11	7.85	5.57	5.36
S-4	934.88	932.3	2.58	10.51	7.93	9.98	7.40	11.43	8.85	8.06
P-5	948.43	946.2	2.23	21.22	18.99	19.59	17.36	21.17	18.94	18.43

DTW BGL Arithmetic Mean of all wells

12.50

TOC - Top of casing

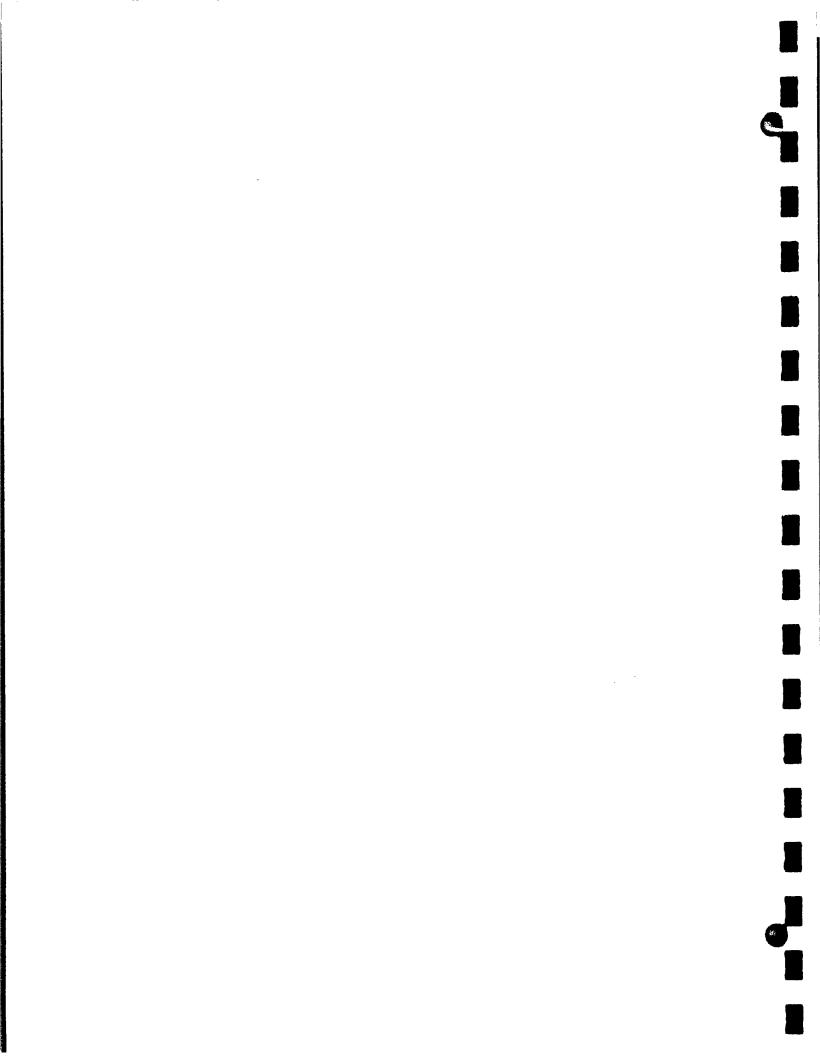
GL - Ground Level

DTW - Depth to water

EKCO\SHALDTW.XLS

BGL - Below ground level

Page 1



Roy F. Weston, Inc. - Gulf Coast Laboratories

DEU Datab No	hav. 04050055	VC	LATILES BY GC/			Report Date: 05	
KFW Batth Nu	<u>ımber: 9405G855</u>	Client: Ekco	Houseware	work	Order: 02994-	002-005-9	<u>Page: 3a</u>
	Cust ID:	R-1	R-2	R-2	R-3	R-3	R-5
Sample Information	RFW#:	007 WATER	008 Water	008 DL	009	009 BL	010
INTOTINACION	Matrix:	WATER 1.0		WATER	WATER	WATER	WATER
	D.F.:		1.0	5.0	1.0	20.0	1.0
	Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
	Toluene-d8	96 %	100 %	104 %	105 %	108 %	98 %
Surrogate	Bromofluorobenzene	102 %	98 %	106 %	101 %	111 %	106 %
Recovery	1,2-Dichloroethane-d4	100 %	98 %	110 %	105 %	112 %	111 %
*********		f)-	:=====f}=	• • • • • • • • • • • • • • • • • • • •	:===########f]		======f1
Chloromethan		2 U	2 U	NA	2 U	NA	2 U
Bromomethane		2 U	2 U	NA	2 U	NA	2 U
Vinyl chlori		1 U	1 U	NA	1 U	NA	 15
Chloroethane		2 U	2 U	NA	2 U	NA	2 U
Methylene Ch	nloride	1 U	1 U	NA	1 U	NA	1 U
Acetone		2 U	5 BU	NA	2 U	NA	2 U
Carbon disul		1 U	1 U	NA	1 U	NA	1 U
1,1-Dichloro		1 U	1	NA	16	NA_	1 U
1,1-Dichloro		1 U	12	NA	E	150	2
	ethene (total)	1 U	E	66	4	NA	17
Chloroform		1 U	1 U	NA	1 U	NA	1 U
1,2-Dichloro	ethane	1 U	1 U	NA	1 U	NA	1 U
2-Butanone		2 U	2 U	NA	2 U	NA	2 U
1,1,1-TrichT		5	1 U	NA	E	-NA-70 ·	1 U
Carbon Tetra	chloride	1 U	1 'U	NA	1 U	NA	1 U
Vinyl acetat	:e	2 U	2 U	NA	2 U	NA	2 U
Bromodich1or	omethane	1 U	1 U	NA	1 U	NA -	1 U
1,2-Dichloro	propane	1 U	1 U	NA	1 U	NA	1 U
cis-1,3-Dich	loropropene	1 U	1 U	NA	1 U	NA	1 U
Trichloroeth		9	21	NA	E	81	5
Dibromochlor	omethane	1 U	1 U	NA	1 U	NA	1 U
1,1,2-Trich1	oroethane	1 U	1 U	NA	1 U	NA	1 U
Benzene		1 U	1 U	NA	1 U	NA	1 U
trans-1.3-Di	chloropropene	1 U	1 U	NA	1 U	NA	1 U
Bromoform		1 U	1 U.,		1 U	NA	1 U
4-Methy1-2-p	entanone	2 U	2 U	·· · · NA · :	2 U	··· NA	2 U
2-Hexanone		2 U	2 U	NA ,	11 2" U !	t NA 😘	2 U
Tetrachloroe	thene	1 U	1 U	NA	1 U	NA	1 U
1,1,2,2-Tetr	achloroethane	1 U	1 U	NA	1 U	NA	1 U
*= Outside o	of EPA CLP QC Timits.						

RFW Batch Number: 9405G855	Client: Ekco He	ouse ware	Work_(Order: 02994-0	002-005-9	Page: 3b
Cust ID:	R-1	R-2	R-2	R-3	R-3	R-5
RFW#:	007	800	008 DL	009	009 DL	010
Toluene	1 0	<u>1</u>	NA	1 U	NA	1 · U ·
Chlorobenzene	1 U	1 U	NA	1 U	NA	1 U
Ethylbenzene	1 U	1 U	NA	1 U	NA	1 U
Styrene	` 1 U	1 U	NA	1 U	NA	1 U
Xylene (total) *= Outside of EPA CLP QC limits.	i Ū	1 Ü	NA	1 U	NA	1 U
· · · · · · · · · · · · · · · · · · ·						

Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS. HSL LIST

R-1 007 WATER 1.0	R-2 008	R-2	R-3	. R-3	R-5
WATER					W-2
ug/L	WATER 1.0 ug/L	008 DL WATER 5.0 ug/L	009 WATER 1.0 ug/L	009 DL WATER 20.0 ug/L	010 WATER 1.0 ug/L
96 % 102 % 100 %	100 % 98 % 98 %	104 % 106 % 110 %	105 % 101 % 105 %	108 % 111 % 112 %	98 % 106 % 111 %
2 U 2 U 1 U 2 U 1 U 2 U 1 U	2 U 2 U 1 U 2 U 1 U 5 B 1 U	NA NA NA NA NA NA NA	2 U 2 U 1 U 2 U 1 U 2 U 1 U	NA NA NA NA NA NA NA	2 U 2 U 15 2 U 1 U 2 U 1 U
1 U 1 U 1 U 2 U 5 1 U	E 1 U 1 U 2 U 1 U 1 U	66 NA NA NA NA NA	4 1 U 1 U 2 U E 1 U	NA NA NA NA 70 NA	2 17 1 U 1 U 2 U 1 U 1 U
2 U 1 U 1 U 1 U 9	2 U 1 U 1 U 1 U 21 1 U	NA NA NA NA NA	2 U 1 U 1 U 1 U E 1 U	NA NA NA NA 81 NA	2 U 1 U 1 U 1 U 5
1 U 1 U 1 U 1 U 2 U 2 U	1 U 1 U 1 U 1 U 2 U 2 U	NA NA NA NA NA NA	1 U 1 U 1 U 1 U 2 U 2 U	NA NA NA NA NA NA	1 U 1 U 1 U 1 U 2 U 2 U 1 U
	102 % 100 % ===================================	102	102 % 98 % 106 % 100 % 98 % 110 %	102	102

RFW Batch Number: 9405G855	Client: Ekco H	ouseware	Work (Order: 02994-0	02-005-9	Page: 3b
Cust ID:	R-1	R-2	R-2	R-3	R-3	R-5
RFW#:	007	800	008 DL	009	009 DL	010
Toluene		<u>1</u>	-NA	1 0	NA	<u>1</u> U
Chlorobenzene	1 U	ע נ	NA	ט נ	NA	j U
Ethylbenzene	1 U	1 U	NA	1 U	NA	1 U
Styrene	1 U	1 U	NA	1 U	NA	1 U
Xylene (total) *= Outside of EPA CLP QC limits.	1 U	1 U	NA	1 U	NA	1 U

.



Roy F. Weston, Inc. - Gulf Coast Laboratories

	VOLATILES BY GC/MS,			05/20/94 11:1	1
RFW Batch Number: 9405G855	Client: Ekco Houseware	Work Order: 02	994-002-005-9	Page: 4	a
					-

<u></u>		OT TOILE.	LIVO	Houseware		MUI	K Older. 0233	4-002-003-9	raye: 4a
	Cust ID:	FB-L-1		TB050294	VBLK		VBLK BS	VBLK BSD	VBLK
Sample	RFW#:	011		012	94GVB164-M	B1	94GVB164-MB1	94GVB164-MB1	94GVB165-MB1
Information	Matrix:	WATER		WATER	WATER		WATER	WATER	WATER
	D.F.:	1.	0	1.0	1.	0	1.0	1.0	1.0
	Units:	ug/L		ug/L	ug/L		ug/L	ug/L	ug/L
	Toluene-d8	104	%	100 %	100	%	104 9	98 X	103 %
Surrogate Br	romofluorobenzene	110	%	107 %	97	%	104 9		
	Dichloroethane-d4	110	%	108 %	99	%	100 %	98 %	105 %
Chloromethane		 2	=f1= U	*=======f 2 U) 2	:=†1 U	2 l	,,	
Bromomethane		2	Ŭ	Ž Ŭ	2	Ŭ	ž i		
Vinyl chloride		ĩ	Ŭ	ĩ Ŭ	ī	Ŭ	i i		
Chloroethane		2	Ŭ	2 U	ż	Ŭ	2 (
Methylene Chloric	le ·	ī	Ū	ĩ Ŭ	ĩ	Ŭ	īì		
Acetone		Ž	Ŭ	4 B	5	•	4 E		
Carbon disulfide		ī	Ū	iŪ	i	U	ii		•
1,1-Dichloroether	ne	ī	Ū	ĪŪ	i	Ŭ	108 9		
1,1-Dichloroethar		i	U	1 U	ì	U	1 (
1,2-Dichloroether	ne (total)	1	U	1 U	1	U	i i	i i i	i ū
Chloroform		1	U	1 U	1	U	i l	i i i	
1,2-Dichloroethan	ie	1	U	1 U	1	U	1 1	i i u	l i ū
2-Butanone		2	U	2 U	2	U	2 (l 2 U	2 U
1,1,1-TrichToroet	hane	1	U	- 1 U	1	U	1 l	1 1 U	1 0
Carbon Tetrachlor	ride	1	U	1 · U	1	U	1 l	1 1 U	i u
Vinyl acetate		2	U	2 U	2	U	2 1	J 2 U	2 U
Bromodichlorometh	nane	1	U	1 U	1	U	1 t	l 1 U	i u
1,2-Dichloropropa	ine	1	U	1 U	1	U	1 l	1 U	1 0
cis-1,3-Dichlorop	propene	1	U	1 U	1	U	1 1	l i u	1 U
Trichloroethene		1	U	1 U	1	U	96 9	5 . 95 x	1 U
Dibromochlorometh	nane	1	U	1 U	1	U	1 (J 1 U	1 U
1,1,2-Trichloroet	hane	1	U	1 U	1	U	1 (J 1 1	1 U
Benzene		1	U	1 U	1	U	108		1 U
trans-1,3-Dichlor	ropropene	1	U	1 U	1	U	1 1	· -	1 0
Bromoform		1	U	1 U	1	U	1 1		
4-Methy1-2-pentar	none	2	U	2 U	2	U	2 l		
2-Hexanone		2	U	2 U	2	U	2 l		
Tetrachloroethene		1	U	1 U	1	U	1 (
1,1,2,2-Tetrachlo	roethane	1	U	1 U	1	U	1 t) 1 U	1 U
*= Outside of EPA	A CLP QC limits.								

RFW Batch Number: 9405G855	Client: Ekco Houseware		Wor	-002-005-9	Page: 4b	
Cust ID:	FB-L-1	TB050294	VBLK	VBLK BS	VBLK BSD	VBLK
RFW#:	011	012	94GVB164-MB1	94GVB164-MB1	94GVB164-MB1	94GVB165-MB1
Toluene	ı u	1 0	1 0	104 %	96 %	
Chlorobenzene	1 U	1 U	1 U	100 %	95 %	1 U
Ethylbenzene	1 U	i U	1 U	1 U	1 U	1 U
Styrene	1 U	1 U	1 U	1 U	1 U	1 U
Xylene (total) *= Outside of EPA CLP QC limits.	1 U	I U	I U	1 0	1 U	1 U

Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS, HSL LIST

Report Date: 05/20/94 11:11

RFW Batch Number: 9405G855 Client: Ekco Houseware Work Order: 02994-002-005-9 Page: 5a

	Cust ID:	VBLK BS		VBLK BSD	÷
Sample	RFW#:	94GVB165-ME	B 1	94GVB165-MB	1
Information	Matrix:	WATER		WATER	
	D.F.:	1.0)	1.0	í
	Units:	ug/L		ug/L	
	Toluene-d8	104	%	99	*
Surrogate	Bromofluorobenzene	108	%		љ %
Recovery	1,2-Dichloroethane-d4	110	%	·	љ %
recovery		110			
Chloromethan		2	Ü		Ü
Bromomethane		- 2	Ŭ		Ŭ
Vinyl chlori		- ī	Ŭ		Ŭ
Chloroethane		- 2	Ŭ	_	Ü
Methylene Ch		້ ເ	Ŭ	_	ij
Acetone		- 1	В	•	B
Carbon disul	Ifide	- 7	Ü		ŭ
1,1-Dichlore		- 118 [*]	%	105	4
1,1-Dichlore		_	ũ		ű
	oethene (total)	- i	Ŭ	•	Ŭ
Chloroform	between (total)	- î	Ŭ	-	ŭ
1,2-Dichlord	nethane	- i	Ŭ	•	บั
2-Butanone		- 2	Ŭ	_	ŭ
1,1,1-Trich	loroethane	- i	Ŭ	-	Ŭ
Carbon Tetra		- i	Ŭ	-	Ŭ
Vinyl acetai		- <u>2</u>	Ŭ	-	Ŭ
Bromodichlor		- <u>ī</u>	Ŭ		Ū
1,2-Dichlore		- <u>i</u>	Ŭ	•	Ŭ
	hloropropene	- <u>i</u>	Ŭ	i	Ū
Trichloroeth		103	%	_	%
Dibromochlor		- 1	ũ		Ü
1,1,2-Trich1		- i	Ũ	_	Ŭ
Benzene		112	%	114	%
	chloropropene		Ü		Ü
Bromoform		- i	Ŭ	· Ī	Ū
4-Methy1-2-p	pentanone	2	Ū	2	Ū
2-Hexanone		- 2	Ū	_	Ū
Tetrachloroe	ethene	- <u>ī</u>	Ŭ		Ū
	rachloroethane	- i	Ū		Ū
	of EPA CLP QC Timits.	_			

BENE DINCONCETE GEOTINITIES. IN TELETIA HELLEGAR IN ROLL. MESICOL ENGT AL L

FACSIMILE TRANSMITTAL Glenloch 2S EAX 610-701-7597



Roy F. Westers Inc. 1 Westen Way West Chester, Pennsylvania 19380-1469 6 610-701-3000 = Fax 610-701-3186

TO:	Pat Cunninghann		nt's Telecopy ne #
FROM: TOTAL I	PAGES: And come break	Origina	tor's Telephone #
W.O. #: COMME		well	data
Providing areas of:	g quality environmental management ar	nd consulting	engineering services for over 40 years in th
Analyt Air Qu Water Hazar	Quality/Wastewater dous, Solid, Radioactive Waste and Safety		Life Sciences Strategic Environmental Management Information Management Construction/Remediation
	54 Office	ces Worldwid	le

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RPW 04.03-806/A-06/M



May 5, 1994

Mr. Dennis Lee Ohio Environmental Protection Agency Northeast District Office 2110 East Aurora Road Twinsburg, Ohio 44087

Dear Mr. Lee:

Enclosed please find Ekco Housewares, Inc. Groundwater Reclamation Project Report.

This report is complete and includes actual analysis from Wadsworth Laboratory.

Sincerely,

EXCO BOUSEWARES, INC.

Jeffrey L. Burman

Process Engineering Manager

Enclosures

JLB/baz

Report #100

cc: Paul Tag, Plant Manager Ekco Housewares, Inc. Massillon, Ohio

> Harold Byer Roy F. Weston, Inc. West Chester, Pennsylvania

Daniel Schiltz Ohio Air Pollution Division Canton, Ohio (Analysis Log - Air Stripper Only) Sally Averill
U.S. Environmental Protection
Agency
Chicago, Illinois

Robert Zollner American Home Products New York, New York

EKCO HOUSEWARES CO.

MR01 W-10" WELL 4-19-94 10:00

WD #: M2397101 LAB #: A4D190023-001

MATRII: WATER

DATE RECEIVED: 4/19/94

DATE SAMPLED:

4/19/94

	_	OF 2 REPORTING			Parish Care Care	00	
PARAMETER	RESULT (ug/L)	LIMIT	METHO	ח	EXTRACTION- ANALYSIS DATE	QC BATCH	
FARACID LAK	(44/11)		WATER!	¥	WANTED THE		
Acrolein	STD	500	USEPA	624	04/20/94	4110117	
Acrylonitrile	MD	500	USEPA	624	04/20/94	4110117	
Benzene	MD	50	USEPA	624	04/20/94	4110117	
Bromodichloromethane	IND	50	USEPA	624	04/20/94	4110117	
Bronoform	MD	50	USEPA	624	04/20/94	4110117	
Bromomethane	ND	100	USEPA	624	04/20/94	4110117	
Carbon tetrachloride	ND	50	USEPA	624	04/20/94	4110117	
Chlorobenzene	MD	50	USEPA		04/20/94	4110117	
Dibromochloromethane	ND	50	USEPA	624	04/20/94	4210117	
Chloroethane	3 1TD	100	USEPA	624	04/20/94	4110117	
2-Chloroethyl vinyl ether	MD	100	USEPA	624	04/20/94	4110117	
Chloroform	MD	50	USEPA	624	04/20/94	4110117	
Chloromethane	ND	100	USEPA	624	04/20/94	4110117	
1,2-Dichlorobenzene	IMD	50	USEPA	624	04/20/94	4110117	
1,3-Dichlorobenzene	ND	50	USEPA	624	04/20/94	4110117	
1,4-Dichlorobenzene	ND	50	USEPA	624	04/20/94	4110117	
1,1-Dichlorosthane	91	50	USEPA	624	04/20/94	4110117	
1,2-Dichlorosthane	IND	50	USEPA	624	04/20/94	4110117	
1,1-Dichloroethene	56	50	USEPA	624	04/20/94	4110117	
1,2-Dichlorostheme, Total	260	50	USEPA	624	04/20/94	4110117	
1,2-Dichloropropane	IKTO	50	USEPA	624	04/20/94	4110117	
cis-1,3-Dichloropropene	ND	50	USEPA	624	04/20/94	4110117	
trans-1,3-Dichloropropene	MD	50	USEPA	624	04/20/94	4110117	
Ethylbenzene	ND	50	USEPA	624	04/20/94	4110117	
SURROGATE RECOVERY	3	ACCEPTABLE	LIMITS				
1,2-Dichloroethane-d4	84	(73 - 1	18)		•		
Toluene-d8	98	{ 92 - 1	09)			į	
Bromoflyorobensene .	. 99	<u>(</u> 85 - 1	08)				

NOTE AS RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LINET

SEN. BY: CONCEPT 610-701-7597 + 7-11-94 :11:16AX : ROY F. RESTON INCH ROY F. RESTON INC. = 4-15

EXCO HOUSEMARES CO.

MR01 W-10* WELL 4-19-94 10:00

WO 8: M2397101 LAD 6: A4D190023-001

DATE RECEIVED:

4/19/94

MATRIX: WATER

----- GC/MS Volutiles -----

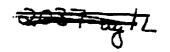
2 OF 2								
PARAMETER	RESULT (ug/L)	REPORTING LIMIT	METHOD	EXTRACTION- ANALYSIS DATE	QC <u>Batch</u>			
Trichlorofluoromethane	MD	50	USEPA 624	04/20/94	4110117			
Methylene chloride	MD	50	USEPA 624	04/20/94	4110117			
1,1,2,2-Tetrachloroethane	ND	50	USEPA 624	04/20/94	4110117			
Tetrachloroethene	MD	50	USEPA 624	04/20/94	4110117			
Toluene	MD	50	USEPA 624	04/20/94	4110117			
1.1.1-Trichlorosthans	1,200	50	USEPA 624	04/20/94	4110117			
1,1,2-Trichloroethane	ЖD	50	USEPA 624	04/20/94	4110117			
Trichlorosthene	430	50	USEFA 624	04/20/94	4110117			
Xylenes, Total	ND	50	USEPA 624	04/20/94	4110117			

SURROGATE RECOVERY	k	ACCEPTABLE LIMITS
1,2-Dichloroethane-d4	84	(73 - 118)
Toluene-ds	98	(92 - 109)
Bromofluorobenzene ·	99	(85 - 108)

NOTE: AN RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LINET





SENT BY: CONCERT GIV-701-7597 - 7-11-94 (11)17AX - ROY F. WESTON, THE ROY F. WESTON, THE

EXCO HOUSEWARES CO.

MRG1 W-10" WELL 4-19-94 10:00

WO #: M2397102

LAB #: A4D190023-001

MATRIX: WATER

DATE SAMPLED: 4/19/94 DATE RECEIVED: 4/19/94

PARAMETER	RESULT (ug/L)	REPORTING LIMIT	HETHOD	EXTRACTION - ANALYSIS DATE	QC <u>BATCH</u>
Vinyl chloride	3.3	1.0	USEPA 624M	04/20/94	4111094

SURROGATE RECOVERY		3	ACCEPTABLE LIMITS
Bromofluorobensene		101	(86 - 115)
1,2-Dichloroethane-64		90	(76 - 114)
Toluene-d8	٠.	98	(88 - 110)

NOTE: AS RECEIVED

*

2040 ug/L

EKCO HOUSEMARES CO.

MR02 SOUTH WELL 4-19-94 10:00

MO #: M2398101 LAB #: A4D190023-002

M4D190023-002

MATRIX: WATER

DATE SAMPLED: 4/19/ DATE RECEIVED: 4/19/:

1 OF 2						
	RESULT	OF 2 REPORTING		EXTRACTION-	QC	
PARAMPTER	(<u>ug/L)</u>	LIMIT	METHOD	ANALYSIS DATE	BAT	
<u> </u>				·		
Acrolein	ND	50	USEPA 624	04/20/94	4110	
Acrylonitrile	XX D	50	USEPA 624	04/20/94	4110	
Bensene	KD	5.0	USEFA 624	04/20/94	4110	
Bromodichloromethane	KID	5.0	USEPA 624	04/20/94	4110	
Bromoform	EID)	5.0	USEPA 624	04/20/94	4110	
Bromomethane	X D	10	USEPA 624	04/20/94	4111	
Carbon tetrachloride	1KD	5.0	USEPA 624	04/20/94	4111	
Chlorobensene	MD	5.0	USEPA 624	04/20/94	411(
Dibromochloromethane	ED	5.0	USEPA 624	04/20/94	4210	
Chloroethane	300	10	USEPA 624	04/20/94	4110	
2-Chloroethyl vinyl ether	ND	10	USEPA 624	04/20/94	4110	
Chloroform	1610	5.0	USEPA 624	04/20/94	411(
Chloromethane	MD	10	USEPA 624	04/20/94	4110	
1,2-Dichlorobensens	XD	5.0	USEPA 624	04/20/94	411(
1,3-Dichlorobenzene	MD	5.0	USEPA 624	04/20/94	4110	
1,4-Dichlorobenzene	ИD	5.0	USEPA 624	04/20/94	4111	
1,1-Dichloroethane	59	5.0	USEPA 624	04/20/94	411(
1,2-Dichloroethane	MD	5.0	USEPA 624	04/20/94	4110	
1,1-Dichlorosthene	6.4	5.0	USEPA 624	04/20/94	4116	
1,2-Dichlorosthene, Total	5. 7	5.0	USEPA 624	04/20/94	411(
1,2-Dichloropropens	36 D	5.0	USEPA 624	04/20/94	411(
cis-1,3-Dichloropropene	RD.	5.0	USEPA 624	04/20/94	4110	
trans-1,3-Dichloropropene) III)	5.0	usepa 624	04/20/94	411(
Rthylbensene	160	5.0	USEPA 624	04/20/94	4110	
SURROGATE RECOVERY	Ì	ACCEPTABLE	LIMITS			
1,2-Dichloroethane-d4	68	(73 - 1				
Toluene-d8	102	(92 - 1				
Aromofluorobensene .	98	(85 - 1	08)			

NOTE: AS RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LIMIT

EXCO HOUSEWARES CO.

MR02 SOUTE WELL 4-19-94 10:00

WO #: M2398101 LAB #: A4D190023-002

LAB #: A4D190023-002 MATRIX: WATER W-1

DATE SAMPLED:

DATE RECEIVED:

4/19/ 4/19/

2 OF 2							
	RESULT	reporting		EXTRACTION-	QC		
PARAMETER	(ug/L)	LIMIT	METHOD	ANALYSIS DATE	<u>BAT</u>		
Trichlorofluoromethane	ND	5.0	USBPA 624	04/20/94	411		
Methylene chloride	197 0	5.0	USEPA 624	04/20/94	411		
1,1,2,2-Tetrachloroethane	RD	5.0	USEPA 624	04/20/94	411		
Tetrachloroethene	ND	5.0	USEPA 624	04/20/94	411		
Toluene	ND	5.0	USEPA 624	04/20/94	411		
1,1,1-Trichlorosthens	23	5.0	USEPA 624	04/20/94	411		
1,1,2-Trichloroethane	ND	5.0	USEPA 624	04/20/94	411		
Trichlorosthens	46	5.0	USEPA 624	04/20/94	411		
Xylenes, Total	MD	5.0	USEPA 624	04/20/94	411		

SURROGATE RECOVERY	Ì	ACCEPTABLE LIMITS
1,2-Dichloroethane-d4	88	(73 - 116)
Toluene-dâ	102	(92 - 109)
Bromofluorobenzene	98	(85 - 108)

NOTE: AS RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LINET

SENT BY: CONCEPT 610-701-7597 : 7-11-94 :11:17AM : ROY F. WESTON, INC. = 8 15

EXCO EGUSERARES CO.

MR02 SOUTH WELL 4-19-94 10:00

13B #: A4D190023-002

MATRIX: WATER

W-1

- - - GC/MS Volatiles - - - -

PARAMETER	RESULT (ug/L)	REPORTINGLIMIT	METHOD	EXTRACTION- AMALYSIS DATE	QC BATCH
Vinyl chloride	MD	1.0	USEPA 624M	04/20/94	4111094

*	ACCEPTABLE LIMITS			
107 101	(86 - 115) (76 - 114) (88 - 110)			
	107			

NO NOT DETECTED AT THE STATED REQUESTING LIBER

EKCO HOUSEWARES CO.

MR03 EFFLUENT FROM AIR STRIPPER 4-19-94 10:00

WO #: M2399101 DATE SAMPLED: 4/19/94
LAB #: A4D190023-003 DATE RECEIVED: 4/19/94

MATRIX: WATER

		/MS Volatiles OF 2			
	result	reporting		EXTRACTION-	ĈC .
PARAMETER	(ug/L)	LIMIT	METHOD	ANALYSIS DATE	BATCH
Acrolein	MD	50	USEPA 624	04/20/94	4110117
Acrylonitrile	IND	50	USEPA 624	04/20/94	4110117
Benzene	190 0	5.0	USBPA 624	04/20/94	4110117
Bromodichloromethane	MD	5.0	USEPA 624	04/20/94	4110117
Bromoform	MD	5.0	USEPA 624	04/20/94	4110117
Bromomethane	KD	10	USBPA 624	04/20/94	4110117
Carbon tetrachloride	19TD	5.0	USEPA 624	04/20/94	4110117
Chlorobenzene	350 0	5.0	USBPA 624	04/20/94	4110117
Dibromochloromethane	350	5.0	USEPA 624	04/20/94	4110117
Chloroethane	NID	10	USEPA 624	04/20/94	4110117
2-Chloroethyl vinyl ether	ND	10	USBPA 624	04/20/94	4110117
Chloroform	ND	5.0	USEPA 624	04/20/94	4110117
Chloromethane	ND	10	USEPA 624	04/20/94	4110117
1,2-Dichlorobenzene	XID	5.0	USEPA 624	04/20/94	4110117
1,3-Dichlorobensene	ND	5.0	USEPA 624	04/20/94	4110117
1,4-Dichlorobenzene	ND	5.0	USBPA 624	04/20/94	4110117
1,1-Dichloroethane	150 0	5.0	USEPA 624	04/20/94	4110117
1,2-Dichloroethane	ND	5.0	USEPA 624	04/20/94	4110117
1,1-Dichloroethene	ND	5.0	USEPA 624	04/20/94	4110117
1,2-Dichloroethene, Total	ND	5.0	USEPA 624	04/20/94	4110117
1,2-Dichloropropane	NIC	5.0	USEPA 624	04/20/94	411011
cis-1,3-Dichloropropene	MD	5.0	USEPA 624	04/20/94	4110117
trans-1,3-Dichloropropene	ND	5.0	USEPA 624	04/20/94	411011
Ethylbenzene	ND	5.0	USEPA 624	04/20/94	4110111
SURROGATE RECOVERY	<u>•</u>	ACCEPTABLE	LIMITS		
1,2-Dichloroethane-d4	89	(73 - 1	-		
Toluene-d8	98	(92 - 1	r 03)		

NOTE: AS RECEIVED

Bromofluorobensene ·

ND NOT DETECTED AT THE STATED REPORTING LIMIT

(.85 - 108)

100

EKCO ECUSEMARES CO.

MR03 EFFLUENT FROM AIR STRIPPER 4-19-94 10:00

WO #: 12399101

DATE SAMPLED:

LAB #: A4D190023-003

DATE RECEIVED:

MATRIX: WATER

• • • • • • • • • • • •	ac/ : 2	MS Volatiles OF 2			• • • •
PARAMETER	RESULT (ug/L)	REPORTING LIMIT	METHOD	EXTRACTION- ANALYSIS DATE	QC <u>Batch</u>
Trichlorofluoromethane	MD	5.0	USEPA 624	04/20/94	411011
Methylene chloride	MD	5.0	USEPA 624	04/20/94	411011
1,1,2,2-Tetrachloroethane	MID	5.0	UBEPA 624	04/20/94	411011
Tetrachloroethene	BTD	5.0	USEPA 624	04/20/94	411011
Toluene	20 0	5.0	USEPA 624	04/20/94	411011
1,1,1-Trichloroethane	SED	5.0	DSEPA 624	04/20/94	411011
1,1,2-Trichloroethane	RD	5.0	USEPA 624	04/20/94	411011
Trichloroethene	KD	5.0	USEPA 624	04/20/94	411011
Ivlenes. Total	ND	5.0	USEPA 624	04/20/94	411011

SURROGATE RECOVERY	ATE RECOVERY + ACC	
1,2-Dichloroethane-de	89	(73 - 118)
Toluene-d8	98	(92 - 109)
Bromofluorobensene	· · 100	(85 - 108)

NO NOT DETECTED AT THE STATED REPORTING LINET

EKCO HOUSEWARES CO.

MR03 EFFLUENT FROM AIR STRIPPER 4-19-94 10:00

MO #: M2399102 DATE SAMPLED: 4/19/94
LAB #: A4D190023-003 DATE RECEIVED: 4/19/94

MATRIX: WATER

PARAMETER	RESULT (ug/L)	reporting Lingt	METHOD	EXTRACTION- ANALYSIS DATE	QC <u>BATCH</u>
Vinyl chloride	ND	1.0	USEPA 624M	04/20/94	4111094

SURROGATE RECOVERY	Ł	ACCEPTABLE LIMITS
Bromofluorobenzene	104	(86 - 115)
1,2-Dichloroethane-d4	95	(76 - 114)
Toluene-d8	· · · 9 7	(88 - 110)

note as received

NO NOT DETECTED AT THE STATED REPORTING LIMIT

RECO MOUSEWARRS CO.

MR04 OUTFALL #001 4-19-94 10:00

WD #: M2400101 LAD #: A4D190023-004 DATE SAMPLED: DATE RECEIVED:

MATRIX: WATER

•	RESULT	OF 2 REPORTING		EXTRACTION-	QC
PARAMETER	(ug/L)	LIMIT	METHOD	ANALYSIS DATE	BATCH
Acrolein	MD	50	USEPA 624	04/20/94	4110117
Acrylonitrile	MID.	50	USEPA 624	04/20/94	4110117
Benzene	MD	5.0	USEPA 624	04/20/94	4110117
Bromodichloromethane	MD	5.0	USEPA 624	04/20/94	4110117
Brancform	MD	5.0	USEPA 624	04/20/94	4110117
Bromomethane	BID	10	USEPA 624	04/20/94	4110117
Carbon tetrachloride	15 0	5.0	USEPA 624	04/20/94	4110117
Chlorobensene	MD	5.0	USEPA 624	04/20/94	4110117
Dibromochloromethane	ND	5.0	USEPA 624	04/20/94	4110117
Chloroethane	BITO	10	USEPA 624	04/20/94	4110117
2-Chloroethyl vinyl ether	MD	10	USEPA 624	04/20/94	4110117
Chloroform	MD	5.0	USEPA 624	04/20/94	4110117
Chloromethane	MD	10	USEPA 624	04/20/94	4110117
1,2-Dichlorobensene	MD	5.0	USEPA 624	04/20/94	4110117
1,3-Dichlorobenzene	MD	5.0	USEPA 624	04/20/94	4110117
1,4-Dichlorobensene	1970	5.0	USEPA 624	04/20/94	4110117
1,1-Dichloroethane	ND	5.0	USEPA 624	04/20/94	4110117
1,2-Dichloroethane	MD	5.0	USEPA 624	04/20/94	4110117
1,1-Dichlorosthese	ND	5.0	USEPA 624	04/20/94	4110117
1,2-Dichlorosthene, Total	ND	5.0	USEPA 624	04/20/94	4110117
1,2-Dichloropropane	MD	5.0	USEPA 624	04/20/94	4110117
cis-1,3-Dichloropropens	MD	5.0	USEPA 624	04/20/94	4110117
trans-1,3-Dichloropropene	ND	5.0	USEPA 624	04/20/94	4110117
Ethylbensene	ND	5.0	USEPA 624	04/20/94	4110117
SURROGATE PROOVERY	ì	ACCEPTABLE	LIMITS		
1,2-Dichloroethane-d4	88	(73 - 1	18)		

(92 - 109)

(-85 - 108)

NOTE: AS RECEIVED

Toluene-d8

NO NOT DETECTED AT THE STATED REPORTING LINET

Bromofluorobensene · · · 101

98

SENT BY: CONCEPT 610-701-7597 : 7-11-94 :11:18AM : ROY F. WESTON, INC- ROY F. WESTON, INC. :#1- 1:

EXCO HOUSEWARES CO.

MR04 OUTFALL #001 4-19-94 10:00

WO #: M2400102 LAB #: A4D190023-004 DATE SAMPLED:

4/19/94

MATRIX: WATER

DATE RECEIVED:

4/19/94

PARAMETER	RESULT (ug/L)	REPORTING LIMIT	METHOD	EXTRACTION- ANALYSIS DATE	QC BATCH
Vinyl chloride	NED	1.0	USEPA 624M	04/20/94	4111094

 SURROGATE RECOVERY
 1
 ACCEPTABLE LIMITS

 Bromofluorobenzene
 100
 (86 - 115)

 1,2-Dichloroethane-d4
 94
 (76 - 114)

 Toluene-d8
 - 99
 (88 - 110)

NOTE AS RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LIBERT

14 ug/L

EKCO ECUSERARES CO.

MR04 OUTFALL #8001 4-19-94 10:00

WO #: M2400101 1.30 #: A4D190023-004

DATE SAMPLED: DATE RECEIVED:

MATRIX: WATER

	2	OF 2			
	RECULT	REPORTING		EXTRACTION-	⊘ C
PARAMETER	(ug/L)	<u> Lucir</u>	METHOD	ANALYSIS DATE	BATCH
Trichlorofluoromethane	36 00	5.0	USEPA 624	04/20/94	4110117
Methylene chloride	SID	5.0	USEPA 624	04/20/94	4110117
1,1,2,2-Tetrachloroethane	MD	5.0	USEPA 624	04/20/94	4110117
Tetrachloroethene	NO.	5.0	USEPA 624	04/20/94	4110117
Toluene	ND	5.0	USEPA 624	04/20/94	4110117
1,1,1-Trichlorosthans	11	5.0	USEPA 624	04/20/94	4110117
1,1,2-Trichloroethane	190 0	5.0	USEPA 624	04/20/94	4110117
Trichloroethene	4.1 3	5.0	USEPA 624	04/20/94	4110117
Xylenes, Total	ND	5.0	USEPA 624	04/20/94	4110117

SURROGATE RECOVERY	1	ACCEPTABLE LIMITS		
1,2-Dichloroethane-d4	88	(73 - 118)		
Toluene-d8	98	(92 - 109)		
Bromofluorobenzene	- 101	(· 85 - 108)		

NOTE AS RECEIVED

NO NOT DETECTED AT THE STATED REPORTING LINET

J ESTEMATED VALUE, (DETECTED), BUT BELOW QUANTITATION LINET.

EKCO HOUSEWARES, INC. - MASSILLON, OHIO SAMPLE ANALYSIS LOG

INFLUENT TO AIR STRIPPER	RATE AT TIME OF	SOUTH WELL W-10" WELL ANALYSIS (Weigh'd Ave	TOTAL. WEIGHT OF VOC'S TO ATMOSPHERE	EPPLUENT FROM AIR STRIPPER SAMPLES	PUMPING RATE AT TIME OF SAMPLES	SAMPLE ANALYSIS (ug/L)	TOTAL GAL. OF WATER TREATED TO DATE	OUTFALL #001 DISHCARGE SAMPLES	SAMPLE TO ANALYS (ug/L)
2.3.93	Som	12062	S. J. Slow	2.3.93	SLOCAD	ND	2002500	 7-3.93	الصلطاء الصلطاء
3233	5306Pm	862.4	5.5 Der.	3.2.93	53000	ND	30,329,3cr	3.2.93	334
41.93	Blocar	0.1911	5.0.Cin	4-1.93	5-KXEW	22mg/	373353r.		Idual .
5-3-93	5410GPM	1004.0	7.016s	5-3-93	540GPM	47 with	25 321,000	 5-3-93	371.00
6-1-93	540GPM	914.0*	10.3 165*	6-1-93	540GPM	NDO	ചാരുന്ന്	6-1-93	41 0 8
7-1-93	535GP/	985.3	11.4165	7-1-93	535GPM	13 ugl	22,046,400	 74-93	63.0
8-9-93	515GP	969.4	10.1165	8-9-93	5/5GP1	NO	28 428 600	 8-9-93	39 0 =
9-2-43	490GPM	1070.G	10.6 lbs	9-2-93	490GP3		16 SED 400	9-2-93	4410
10-1-93	464 GP1	1018.6	9 165	10-1-93	4645PM	()	19,679,500	 10-1-93	41,00
11-3-93	494 GPM	14a2*	8.4165	11-3-93	1 . I		19 044 200	11-3-93	63us
12-1-93	300GPM	1263	4.6 165	12-1-93	300GPA	NB	12,173,200	12-1-93	26 vs 7
1-13-94	590GPM	823	11.8165	13-13-94	590GP/	QuelL	37.560.000	1-13-94	69 mg >
2-14-94	440GPM	1238	10.1165	2.14.94	440GPM	3.01	22.841 800	2-14-94	47.60
3-3-94	540GPM	1081	13.0 lbs	3-3-94	540gpm	άN	17642400	3.3.94	49 115
4-19-94	535GPM	1170	14.0 lbs	4-19-94	535GPM		2G. 097.200	4-19-94	15 uc.
									15 ug 7.55 (o)
									9
									7

#Estimated (Pump meter #2 not working)

*Estimated (Southwell meter not working)

Weights nug for with well only

SENT BY CONCE-

\(\(\);=15 ic

SOUTH WELL SAMPLES	WATER LEVEL AT TIME OF SAMPLE	PUMPING RATE AT TIME OF SAMPLE	TOTAL GAL. OF WATER TREATED TO DATE	SAMPLE ANALYSIS (ug/L)	W-10" Weill Samples	WATER LEVEL AT TIME OF SAMPLE	PUMPING RATE AT TIME OF SAMPLE	TOTAL GAL. OF WATER TREATED TO DATE	SAMPLE ANALYS IS (ug/L)
2.3.93	120	245 GPM	903 ETGP	18341	2.383	57'	315 GPm	11389800	Z16SmJr
37.43	155,	2:30(5Pm	2005,000	135m/L	3-2.93	35	300 GPM	//7 4/3 00	Marash
<u> 1-143</u>	122,	5.5K. 2011	5332 tan	Prose	41.93	37'	310GR11	15tactor	195462
<u>5-3-93</u>	1301	235 GPM	10,93800	136 mg/L	5-3-93	851	305 GPM	14 383,000	1673
6-1-43	115'	230GPM	9 408 000	NB	6-1-93	68'		# 12,623.600	
7-1-93	121'	235 GPM	9,553,000	124 ug/L	7-1-93	יגל.		12 493 900	() (
8-9-93	105'	SIDGEM	11 727 000	132 wall	8-9-93	85′		IG. 701,000	
7-2-93	1054	170 GPM	ര ക്ഷേത	140 m/L	9-2-93	80'		10 205,000	(
0-1-93	1001	175 GPM	7, 505 m	136 ucl	10-1-93	80'		12.194.500	3
1-3-93	100'	* Saxte	2000000	* 0	11-3-93	751		11 044 200	7.
2-1-93	951		operation	ıl	12-1-93	751		מגנדוגו	1263 w/L
-13-94	90'	300 GPM*	19,000,000	136 walL	1-13-94	80'		18 560 500	7.
1-14-94	90'	145GPM	961200	138 w. L	2-14-94	72'		13,229,800	<u> </u>
3-3-94	80'	245 GPM	10,584,000	7	3-3-94	50'		7,058,400	
1-19-94	125'		14,102,000	141 101	4-19-94	80'	A .	11,995 200	Z.
				9.					0
									

* Estimated (Pump meter 2 not working)

t meter not working (estimated)

* functioning

RFW Batch Number: 9203G477	Client: Ekco	Houseware	Wor	Page: 1b			
Cust ID:	1-7-GW	1-7-GW	1-8-GW	D-4-30-GW	D-4-30-GW	D-4-30-GW	
RFW#:	001	01 001 DL 002		003	003 DL	003 DL	
Toluene	1 U	NA	1 U	1 U	NA	NA	
Chlorobenzene	. 1 U	NA	1 U	1 U	NA	NA	
Ethy1benzene	. 1 U	NA	1 U	1 U	NA	NA	
Styrene	. 1 U	NA	1 U	1 U	NA	NA	
Xylene (total)	์ 1 บ	NA	1 U	1 บ	NA	NA	
*= Outside of EPA CLP QC limits.	•						

in the second

- (100 to 100 to

Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS. HSL LIST

RFW Batch Number: 9203G477		VOLATILES BY GC/MS, Client: Ekco Houseware					<u>k Order: 299</u>	4-0	керогі vate 2-03-0000	: 04/20/9 Pa	04/20/92 10:51 Page: 1a		
Cust ID:		1-7-GW		1-7-GW		1-8-6M		D-4-30-GW		D-4-30-GW	D-4-30-GW		
Sample	RFW#:	001		001 DL		002		003		003 DL	00	3 DL	ì
Information Matrix:		WATER		WATER	_	WATER	4	WATER		WATER			
D.F.	D.F.:	1.00		20.0		1.00		1.00		20.0		WATER 400	
	Units:	ug/l				ug/L		ug/L		ug/L		ug/L	
	Toluene-d8	99	*	94	*	104	*	106	×	99	% 1	01	%
Surrogate Bromof	luorobenzene	102	%	94	%	99	×		ž		-	00	*
Recovery 1,2-Dichl	oroethane-d4	83	%	89	%	77	%	83	%	98	%	94	×
Chloromethane		2	U	NA		2	† 1 ·	2	:T1= U	NA	,,	A	f1
		2	Ŭ	NA		2	Ŭ		Ŭ	NA		A	
Vinyl Chloride		ī	Ū	NA		ĭ	Ū	_	Ŭ	NA		A	
		7.2	•	NA		Ž	Ŭ	_	Ŭ	NA		A	
Chloroethane Methylene Chloride		3	u	NA		3	-	Ă	Ŭ	NA		A	
ACETONE		2	U	NA		2	U	-	U	NA		Α	
Carbon Disulfide		1	U	NA		ĭ	Ŭ	ī	Ū	NA		A	
1,1-Dichloroethene		. 5.	1 .	NA		ī	Ū	_	Ě	*310		iA .	
1.1-Dichloroethane			E	»14Q		ī	Ū		Ē	A10		IA .	
1,2-Dichloroethene (t	otal)	1	Ü	NA		i	Ū		Ē	730		iA .	
Chlamafaum		i	Ū	NA		ī	Ŭ	91		NA		IA	
1,2-Dichloroethane		ĩ	Ū	NA		ī	Ū	_	U	NA		iA	
2-Rutanone		2	U	NA		2	Ŭ	Ž	Ŭ	NA ·		A	
1,1,1-Trichloroethane		711		NA		ī	Ŭ	_	Ĕ			700	•
carbon recracilloride		1	U	NA		ī	Ū	1	Ũ	KA		IA	
vinyi Acetate		2	U	NA		2	Ū	Ž	Ŭ	NA		ΪA	
Bromodichloromethane		Ī	U	NA		Ī	Ū	Ī	Ŭ	NA		iA	
4 A B1.L1		Ī	IJ	NA		Ĭ	Ŭ	Ī	Ū	NA		iA	
cis-1,3-Dichloroprope	ne	1	U	NA		Ĭ	Ū	Ī	Ü	NA		iA .	
Trichloroethene		914		NA		ī	•	-	Ē		E		E
Dibromochloromethane		i	U	NA		i	U	1	Ũ	NA	_	iA	_
1,1,2-Trichloroethane		ì	Ü	NA		ī	Ŭ	ī	Ū	NA		IA .	
Benzene	•	ī	Ū	NA		i	Ŭ	ī	Ū	NA		IA .	
Trans-1,3-Dichloropro	pene	i	Ū	NA		i	Ū	ī	Ŭ	NA		IA .	
		i	Ŭ	NA		ī	Ŭ	ĺ	Ŭ	NA		IA	
4-Methyl-2-pentanone		2	Ū	NA		Ž	Ŭ	Ž	Ŭ	NA		iA	
Z-Hexanone		2	Ŭ	NA		2	Ŭ	Ž	Ŭ	NA		IA .	
Tetrachloroethene		ĩ	Ū	NA		ī	Ŭ	19 74	· •	NA		ÍΑ	
1.1.2.2-Tetrachloroe Outside of EPA CL	thana	ī	Ù	NA		ī	Ū	ĭ	u	NA		ÍΛ	

RFW Batch Number: 9203G478 Client; Ekco Houseware Work Order: 2994-02-03-0000 Page: 8b

RFW#: 92GVE092-MB1

Chlorobenzene	120 %	
1,1,1,2-Tetrachloroethane	0.2 U	
Ethylbenzene	0.1 U	
Styrene	0.2 U	
p-Xylene	0.2 U	
1,1,1,2-retrachioroethane Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane	0.2 U	
o-Xylene	0.3 U	
Bromobenzene	0.1 U	
1,2,3-Trichloropropane	0.4 U	
Isopropylbenzene	0.2 U	
n-Propylbenzene	0.2 U	
2-Chlorotoluene	0.2 U	
Isopropylbenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene 1,3,5-Trimethylbenzene	0.2 U	
1.3.5-Trimethylbenzene	0.1 Ŭ	
tert-Butvlbenzene	0.2 U	
1.2.4-Trimethylbenzene	0.3 U	
tert-Butylbenzene 1,2,4-Trimethylbenzene sec-Butylbenzene p-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene	0.2 U	
n-Isopropyltoluene	0.2 U	
1.3-Dichlorobenzene	0.1 U	
1.4-Dichlorobenzene	0.3 Ŭ	
1.2-Dichlorobenzene	0.1 U	
n-But v 1 henzene	0.1 ΰ	
n-Butylbenzene 1,2-Dibromo-3-Chloropropane	0.5 U	
1,2,4-Trichlorobenzene	0.2 U	
Hexachlorobutadiene	0.4 U	
Naphthalene 1,2,3-Trichlorobenzene	0.2 U	
1.2.3-Trichlorobenzene	0.1 U	
cis-1,3-Dichloropropene	0.1 U	
trans-1,3-Dichloropropene	0.3 U	
Acetone		
Carbon Disulfide	0.1 U	
2-Butanone	0.3 Ŭ	
Vinyl Acetate	0.3 Ü	
2-Butanone Vinyl Acetate 4-Methyl-2-pentanone	0.4 U	
2-Hexanone	0.5 U	
2-Chloroethylvinylether	0.2 U	
*= Outside of EPA CLP QC limits.	·	

RFW Batch Number: 9203G4	78	V Client:	OLAT Ekco	ILĖS BY GC/MS, Houseware	METHOD		Report Date 2994-02-03-0000	: 04/15/92 11:21 Page: 8a
	Cust ID:	VBLK BS						
Sample Sample	RFW#:	92GVE092-H	R1					
Information	Matrix:	WATER	-					
	D.F.:	1.0	0					
	Units:	UG/L						
Bromofluo	robenzene		%			<u> </u>	f]	<u> </u>
Dichlorodifluoromethane_	******	0.3						
Chloromethane	_	- 0.2						
Bromomethane		0.1						
Bromomethane Vinyl Chloride		- 0.2						
Chloroethane		0.1						
Chloroethane Methylene Chloride Trichlorofluoromethane	·	0.6	Ŭ					
Trichlorofluoromethane		0.1	Ū					
1,1-Dichloroethene		162 *						
1.1-Dichloroethane		0.1						
is-1,2-Dichloroethene_		0.2	Ü					
2,2-Dichloropropane		0.1	U					
trans-1,2-Dichloroethene		0.1						
Chloroform -		0.2	U					
Bromochloromethane		0.2	U					
l.2-Dichloroethane		0.2	U					
l.l.l-Trichloroethane		0.2	U					
Carbon Tetrachloride		0.1	U					
l.l-Dichloropropene		0.1	U					
Bromodichloromethane		0.2	U					
Dibromomethane		0.2	U					
1,2-Dichloropropane		0.2						
Trichloroethene		200 *	%					
Dibromochloromethane		0.2	U					
1.2-Dibromoethane		0.2	U					
1,1,2-Trichloroethane		0.3	U					
Benzene		118	%				•	
1,3-Dichloropropane		0.2						
Bromoform		_ 0.2						
Tetrachloroethene		0.1						
1,1,2,2-Tetrachloroethan	9	0.2						
Toluene		111	%					
- Outside of EPA CLP QC	limits.	-						<u> </u>

RFW Batch Number: 9202	G419	•	V	OLATILES BY	GC	C/MS, HSL LI	ST Wor	k Order: 2994-1	Report Date 02-02-0000	: (D2/25/92 l <u>Page:</u>	4:49
	Cust ID:	Trip Blank	•	Trip Blank	;	FB-L-3	:	FB-L-3	L-1		L-	-1
Sample	RF₩#:	001		002		003		003 DL	004		004 D	
Information	Matrix:	WATER		WATER		WATER		WATER	WATER		WATER	
	D.F.:	1.0	10	1.0	0	1.0	0	4.00	1.00)		.00
	Units:	ug/L	•	ug/L		ug/L		ug/L	ug/L		ug/	/ L
	Toluene-d8	100	*	103	×		*	102 %		7	104	*
	uorobenzene	97	*	101	*		*	107 %	•	%	109	*
Recovery 1,2-Dichlo	roethane-d4	. 99	% £1	101	% f1	106 }	% _f1	105 %		% .f1	112	% f1
		_	Ü	2	U	2	Ü	. NA	2	Ü	NA	,
Bromomethane		2	U	2	U	2	U	NA	2	Ų	NA_	_
Vinyl Chloride		1	U	1	U	1	U	NA		E	712	1
Chloroethane		- 2	U	2	U	2	U	NA	_	U	NA	
Chloroethane Methylene Chloride	······································	- <u>1</u>		0.9	J	5		NA	1	Y	. NA	
Acetone		- 6		2	U	2	U	NA	2		NA	
		- i	บ	1	U	1	U	· NA	1	U	NA	
			U	1	U	1	U	NA	4 8 ,	١	NA	
1 1 D1 - 1 1 Al		1	U	1	U	1	U	NA		Ē	159	_
1,1-Dichloroethane (to	tal)	1	U	1	u	1	U	NA	_	E	5 9 (י ס
Chloroform		_ 1	U	1	U	2		NA	1	U	NA	
1,2-Dichloroethane		1	U	1	U	•	U	NA	1	U	· NA	
2-Butanone		2	U	2	U	2	U	NA	2	Ų	NA	
1.1.1-Trichloroethane		1	U	1	U	1	U	NA	144.5	4	NA	
Carbon Tetrachloride_		1	บ	1	U	1	U	NA	1	U	KA	
Yinvi Acetate		2	U	Z	IJ	2	U	NA	2	U	NA	
Bromodichloromethane_		- 1	U	1	U	1	U	NA.	1	IJ	NA	
1.2-Dichloropropane			U	1	U	1	U	NA	1	U	NA	
cis-1,3-Dichloropropen	A	- <u>ī</u>	Ü	1	U	1	U	NA	1	U	NA	
Trichloroethene	~	- i	ŭ	Ĭ	Ü	_	E	26		£		E
Dibromochloromethane_		- j	ū	ì	Ū	1	U	NA	1	U	NA	
1,1,2-Trichloroethane		- j	Ū	i	Ū	1	Ü	NA	1	U	NA	
Benzene		- i	Ũ	Ĭ	Ū	Ĩ	Ü	NA	1	U	NA	
Trans-1,3-Dichloroprop	ene	- i	Ū	ī	Ū	Ī	Ü	NA .	1	U	NA	
Bromoform		1	ŭ	i	Ŭ	_	Ũ	NA NA	1	U	NA	
4-Hethyl-2-pentanone_		- ;	ŭ	ž	Ŭ	2	ŭ	NA	2	Ū	NA	
2-Hexanone		- 2	ŭ	ž	Ũ	Ž	Ŭ	NA	2	U	NA	
Tetrachloroethene		- ī	ŭ	ĭ	Ũ	_	Ü	NA	1	U	NA	
1,1,2,2-Tetrachloroeth	ane	- i	ŭ	i	Ū	•	Ū	NA .	1	U	NA	
*- Outside of EPA CLP	<u>-</u>		•	•	_	-	-	- - •				

Bat water 920	ien	7.15		FB-L-3	32-CE	Park Jane
Cust ID: Tri	p 81 ank T 001	rip Blank 002	FB-L-3 003	003 DL	004	004 DL
Toluene Chlorobenzene Ethylbenzene Styrene Xylene (total) *** Outside of FPA CLP OC limits.	1 U 1 U 1 U 1 U	1 U 1 U 1 U 1 U	0.9 J 1 U 1 U 1 U	NA NA NA NA	1 U 1 U 1 U 1 U	NA NA KA NA NA

RFW Batch Number: 9109G724	Client:	Ekc	o Houseware			Wor	k Order: 29	94-	02-03-0000		Page: 5b
Cust ID:	VBLK BS		VBLK		VBLK BS		VBLK		VBLK BS		
RFW#:	916VC313-P	IB1	916VB342-MB	1	91GVB342-M	B1	91 GVB 343-M	B 1	91GVB343-MB	1	
Toluene	105	*	5	U	92	%	. 5	U	98	7	
Chlorobenzene	108	×	5	U	94	%	5	Ū	0.4	%	
Ethylbenzene	5	U	5	U	5	U	5	Ü		Ü	
Styrene	5	U	5	U	5	U	5	U		Ū	
Xylene (total)	5	U	5	U	5	U	5	U	5	Ū	
2-Chloroethylvinylether *= Outside of EPA CLP QC limits.	10	U	10	U	10	U	10	U	10	U	

Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS, HSL LIST
Client: Ekco Houseware Work Order: IST Report Date: 10/10/91 13:07 Work Order: 2994-02-03-0000 Page: 5a

RFW Batch	Number: 91096724	<u>Client:</u>	Ekc	o Houseware			<u>Wor</u>	k Order: 2994	-02-03-0000		Page: 5a
	Cust ID:	VBLK BS		VB LK		VBLK BS		VBLK	VBLK BS		
Sample Information		91GVC313-M SOIL	B 1	91GVB342-MB Water	1	91GVB342-M WATER	B1	91GVB343-MB1 WATER	91GVB343-M WATER	Bl	
200000000000000000000000000000000000000	D.F.:	1.0	0	1.00)	1.0	0	1.00	1.0	0	
	Units:	ug/K		ug/L		ug/L		ug/L	ug/L		
	Toluene-d8	96	%		X	101 107	% %	96 % 107 %		%	
Surrogate	Bromofluorobenzene	99 105	% %		% %	107	76 %	107 %		* %	
Recovery	1,2-Dichloroethane-d4					1V7 	-f1		112	F] ====== f]
Chlorometh			Ū		U	10	Ü	10 U	•	Ü	,,,
	ne	10	Ū		Ū	10	Ũ	10 U		Ū	
Vinyl Chlo	ne ride	10	Ū		Ü	10	U	10 U		U	
Chloroetha	ne	10	U	10	U	10	U	10 U	10	U	
Methylene (ne Chloride	25	В	5	U	5	U	5 U	15		
Acetone		31		17		10	U	10 U	10	U	
Carbon Dis	ulfide roethene	⁻ 5	U		U	5	U	5 U		U	
1,1-Dichlo	roethene	101	%	_	U	85	%	5 U		%	
			U		U	5	U	5 U		U	
1,2-Dichlo	roethene (total)	5	U	_	U	5	U	5 U		U	
Chloroform		5	U	•	U	5	U	5 U	•	U	
1,2-Dichloi	roethane	_ 5	U	_	U	5	U	5 U	-	U	
2-Butanone		10	U		U	10	U	10 U		U	
1,1,1-Tricl	hloroethane	_ 5	U	-	U	11		5 U	-	U	
Carbon Teti	rachloride	_ 5	U	_	U	5	U	5 U	•	U	
Vinyl Aceta	ate	10	U		U	10	U	10 U		U	
Bromodichle	oromethane	5	U		U	5	U	5 U	_	U	
1,2-Dichlo	ropropane chloropropene thene	5 5	U		U	5	U	5 U	•	U	
V10 210 011	cirror opi opene		U		U	5	U	5 U	_	U	
11 161110106	thene		%		U	246 *	~	5 U		%	
Dibromochle	oromethane	5	U		U	5	U	5 U		U	
1,1,2-Tricl	oromethane hloroethane	5	U		U	5	U	5 U	_	U	
Benzene	Dichloropropene	118_	%		U	96_	%	5 U		%	
			U		U	5	U	5 U	•	U	
Bromoform		5	U		U	. 5	U	5 U	_	U	
4-Methyl-2	-pentanone	10	Ü		Ü	10	Ü	10 U	• •	U	
Z-Hexanone		: 10	Ü		Ü	10	U	10 U		U	
Tetrachlor	oethene	5	Ü		U	5	Ü	5 U	_	Ü	
	trachloroethane	5	U	5	U	5	U	5 U	5	U	
== Uutside	of EPA CLP QC Timits.										

RFW Batch Number: 9109G724	Client: Ek	co Houseware	Wor	<u>k Order: 2994-</u>	02-03-0000	Page: 4b
Cust ID:	SB07-FB	VBLK	VBLK BS	VBLK	VBLK BS	VBLK
RFW#:	013	91GVB338-MB1	91GVB338-MB1	91GVB339-MB1	91GVB339-MB1	91GVC313-MB1
Toluene Chlorobenzene Ethylbenzene Styrene Xylene (total) 2-Chloroethylvinylether *= Outside of EPA CLP QC limits.	5 U 5 U 5 U 5 U 10 U	5 U 5 U 5 U 5 U 10 U	82 % 93 % 5 U 5 U 10 U	5 U 5 U 5 U 5 U 10 U	85 % 93 % 5 U 5 U 5 U 10 U	5 U 5 U 5 U 5 U 10 U

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Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS, HSL LIST

Report Date: 10/10/91 13:07 Work Order: 2994-02-03-0000 RFW Batch Number: 91096724 Client: Ekco Houseware Page: 4a

Krw Batch Number: 910	730/67	Clianti	EKC	<u>o nouseware</u>		MOL	k order: 2994-	-02-03-0000	Page: 4a
	Cust ID:	SB 07-F	В	VBLK	VBLK BS		VBLK	VBLK BS	VBLK
Sample .	RFW#:	01	3	916VB338-MB1	91GVB338-M	IB1	91GVB339-MB1	91GYB339-MB1	91GVC313-MB1
Information	Matrix:	WATER		WATER	WATER		WATER	WATER	SOIL
	D.F.:	1.0		1.00	1.0	00	1.00	1.00	1.00
	Units:	ug/	L	ug/L	ug/L	-	ug/L	ug/L	ug/Kg
	Toluene-d8	102	*	96 5		*	98 %		97 %
	luorobenzene	102	%	95 5		%	103 %	110 %	
Recovery 1,2-Dichl	oroethane-d4	99	% £1	98 2	112 	% f1	95 %	106 % 1f	
Chloromethane		10	Ü	10 t	10	Ü	10 U	10 Ú	10 Ú
Bromomethane		10		10 l	• • •	U	10 U	10 U	10 U
Vinyl Chloride		10		10 (• • •	U	10 U	10 U	10 U
Chloroethane		10		10 l		U	10 U	10 U	10 U
Chloroethane Methylene Chloride		5	U	5 l			7	5 U	10
Acetone		10		10 l		U	12	32 B	10 U
Carbon Disulfide		5	_	5 l	•	U	5 U		5 U
1,1-Dichloroethene		5	U	5 {		%	5 U	• • • • • • • • • • • • • • • • • • • •	_ :
1,1-Dichloroethane		5	U	5 t	•	U	5 U		5 U
1,2-Dichloroethene (t	total)	5	U	5 t	•	U	5 U	5 U	5 U
Chloroform		5	U	5 t	•	U	5 U		5 U
1,2-Dichloroethane		5	U	5 t	5	U	5 U		5 U
2-Butanone		10	_	10 (U	10 U		10 U
1,1,1-Trichloroethane	9	5	U	5· l	•	U	5 U	5 U	5 U
Carbon letrachioride_		5	_	5 l	•	U	5 U	• -	5 U
Vinyl Acetate		10	U	10 l		U	10 U	10 U	10 U
Bromodichloromethane_		5	U	5 l	_	U	5 U	• -	5 U
1,2-Dichloropropane_cis-1,3-Dichloroprope		5	U	5 t	,	U	5 U		
cis-1,3-Dichloroprope	ne	5	U	5 l	•	U	5 U		
Trichloroethene		5	U	5 t		%	5 U	~	
Dibromochloromethane_		5	U	5 (U	5 U		5 ป
1,1,2-Trichloroethane)	5	U	5 (•	U	5 U		_
Benzene		5		5 (%	5 U		• -
Trans-1,3-Dichloropro	pene	5	U	<u>5</u> (•	U	5 U	• -	
Bromoform		5	U	5 l	_	U	5 U		5 U
4-Methyl-2-pentanone		10		10 (U	10 U		
2-Hexanone		: 10		10		U	10 U		
Tetrachloroethene		5		5 t	_	U	5 U	• -	• -
1,1,2,2-Tetrachloroet		5	U	5 (J 5	U	5 U	5 U	5 U
*= Outside of EPA CLF	QC limits.								

RFW Batch Number: 9109G724 Cust ID:	<u>.Client:</u> SB07-04-06	3	SB07-00-02		R-2-GW	Order: 2994-0 R-2-GW	1-5-6W	Page: 1-6-GW
RFW#:	008 MSD)	009		010	010 DL	011	012
Toluene	94	*	6	U	5 U	NĀ	5 U	5
Chlorobenzene	88	%	6	U	5 Ū	NA	Š Ū	5
Ethylbenzene	6	U	6	U	5 Ū	NA	5 Ū	5
Styrene	6	U	6	U	5 Ū	NA	5 Ü	5
Xylene (total)	6	U	6	U	5 Ū	NA .	5 Ü	5
2-Chloroethylvinylether	11	U	12	U	10 Ū	NA	10 Ū	10
*- Outside of EPA CLP QC limits.	•		되고(-)	1377		118	- H

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Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS, HSL LIST Report Date: 10/10/91 13:07
Client: Ekco Houseware Work Order: 2994-02-03-0000 Page: 3a DEM Ratch Number: 01006724

RFW Batch Number: 91	096724	Client:	Ekc	o Houseware			Wor	<u>k Order: 2994-02</u>	<u>-03-0000</u>		Page:	<u> 3a</u>
	Cust ID:	SB07-04-06	5	SB07-00-02		R-2-GW	1	R-2-GW	I-5-GW		1-6-GW	1
Sample	RFW#:	008 MSI)	009	1	010	t	010 DL	011		012	
Information	Matrix:	SOIL		SOIL		WATER		WATER	WATER		WATER	
	D.F.:	1.0		1.0		1.0		10.0	1.0	0	1.0	10
	Units:	ug/l	(g	ug/K	g	ug/L		ug/L	ug/L		u g/L	•
	Toluene-d8	107	*	107	*	100	%	105 %	101	%	101	7%
	fluorobenzene	88	%	85	X	99	×	109 %	101	%	104	%
Recovery 1,2-D1ch	loroethane-d4	98	% f1	101	% _f1	94	% =f1	113 %	94	% _f1-	97	% f1
Chloromethane		11	Ü	12	Ü	10	Ü	NA	10	Ü	10	Ü
			U	12	U	10	U	NA	10	U	10	U
Vinyl Chloride		11	U	12	U	130		NA	10	U	10	U
Chloroethane		11	U	12	U	10	U	NA	10	U	10	U
Chloroethane Methylene Chloride		33	В	47	В	5	U	NA	5	U	5	U
Acetone		100		37	P	46	В	NA	10	U	10	U
Carbon Disulfide		6	U	6	U	5	U	NA	5	U	5	U
1,1-Dichloroethene_		70	%	6	U	51		NA	5	U	5	U
1,1-Dichloroethane		6	U	6	U		E	240	48	•	5	U
1,2-Dichloroethene (total)	43		210				80 180	5	U	5	U
			U	6	U	5	U	NA	5	U	5	U
Chloroform 1,2-Dichloroethane		6	U	6	U	5	U	NA	5	U	5	U
A D. A		• • •	U	12	U	10	U	NA	10	U	10	U
1,1,1-Trichloroethan	e	6	U	6	U	16		NA	5	U	5	U
Carbon letrachloride		6	U	6	U	5	U	NA	5	U	5	U
Vinyi Acetate		11	U	12	U	10	U	NA	10	U	10	U
Bromodichloromethane		b	U	6	U	5	U	NA	5	U	5	U
1,2-Dichloropropane_			U	6	U	5	U	NA	5	U	5	U
1,2-Dichloropropane_cis-1,3-Dichloroprop	ene	6	U	6	U	5	U	NA	5	IJ	5	IJ
irichioroethene		99	%	210			E	160 760	5	U	5	U
Dibromochloromethane		6	U	6	U	5	U	NA .	5	U	5	U
1,1,2-Trichloroethan	e	6	U	6	U	5	U	NA	5	U	5	U
Benzene		96	%	6	U	5	U	NA	5	U	5	U
Trans-1,3-Dichloropr	opene	6	U	6	U	5	U	NA	5	U	5	U
D		_	U	6	U	5	U	NA	5	U	5	U
4-Methyl-2-pentanone		11	U	12	U	10	U	NA	10	U	10	U
2-Hexanone		11	U	12	U	10	U	NA	10	U	10	U
Tetrachloroethene		, 6	U	6	U	5	U	NA	5	U	5	U
1,1,2,2-Tetrachloroe	thane	6	U	6	U	5	U	NA	5	U	5	U
*= Outside of EPA CL	P QC limits.	-										

RFW Batch Number: 9109G724		Client:	Ekc	co Houseware			Wor	rk Order: 29	94-	-02-03-0000		Page:	2b
Cust	ID:	SB08-02-04		SB08-06-08		SB08-10-12		SB08-10-12 Dup		SB07-04-06		SB07-04-06	
RF	W#:	004		005		006	i	007		800		008 MS	•
Toluene		6	U	6	U	6	U	6	U	6	U	93	*
Chlorobenzene		_ 6	U	6	U	6	U	6	U	6	U	86	%
Ethy1benzene		6	U	6	U	6	U	6	U	6	U	6	U
Styrene		6	U	6	U	6	U	6	U	6	U	6	Ū
Xylene (total)		- 6	U	6	U	6	U	6	U	6	Ū	6	Ĭ
2-Chloroethylvinylether *= Outside of EPA CLP QC limit	•	12	Ū	12	Ū	11	Ū	12	Ū	11	Ũ	11	Ŭ
occordo or cirroti do timio	••	<i>}</i> '	b	V I	b	N	(i)			111			

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Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS, HSL LIST Report Date: 10/10/91 13:07
Client: Ekco Houseware Work Order: 2994-02-03-0000 Page: 2a RFW Batch Number: 9109G724

	Cust ID.							
	cast in:	SB08-02-04	ļ	SB08-06-08	SB08-10-12	SB08-10-12 Dup	SB07-04-06	SB07-04-06
Sample	RFW#:	004		005	006	007	008	008 MS
Information	Matrix:	SOIL	-	SOIL	SOIL	SOIL	SOIL	SOIL
	D.F.:	1.0	00	1.00	1.00	1.00	1.00	1.00
	Units:	ug/l	-	ug/Kg	ug/Kg	ug/Kg	ug/Kg	ug/Kg
		-3/ .	. 3	-3/3	-3/3	-5/3	-3/113	49/119
	Toluene-d8	97	*	104 %	101 9	99 %	103 %	109 %
Surrogate Bromoflu	ıorobenzene	91	%	90 %				85 %
Recovery 1,2-Dichlor		100	%	100 %				98 %
			•=f1	f				======f
Chloromethane		12	U	12 U	11 (12 U	11 U	11 Ù
Bromomethane		12	U	12 U	11 U		11 U	11 U
Vinyl Chloride		12	U	12 U	11 (l 12 U	11 U	11 U
Chloroethane		12	U	12 U	11 U	12 U	11 U	11 U
Methylene Chloride	,	18	₿.		20 E	14 B	30 B	36 B
Acetone		64	N,) 6 4 () 100 ·	் 100 ƙ	93 π	110
Carbon Disulfide		6	U	6 U	6 t	1 6 U	6 U	6 U
1,1-Dichloroethene		6	U	6 U	6 t	1 6 U	6 U	71 %
1,1-Dichloroethane		6	U	6 U	6 L	1 6 U	6 U	6 Ü
1,2-Dichloroethene (tot	al)	6	U	6 U	6 l	l 6 U	31	53
Chloroform		6	U	6 U	6 L	1 6 U	6 U	6 U
1,2-Dichloroethane		6	U	6 U	6 L	l 6 U	6 U	6 U
?-Butanone		12	U	12 U	11 4	12 U	11 U	11 0
1,1,1-Trichloroethane		6	U	6. U	6 U	1 6 U	6 U	6 U
Carbon Tetrachloride		6	U	6 U	6 l	l 6 U	6 Ü	6 Ū
Vinyl Acetate		12	U	12 U	11 U	12 U	11 Ū	li Ū
Bromodichloromethane		6	U	6 U	6 U		6 U	6 0
1,2-Dichloropropane		6	U	6 U	6 L	1 6 U	6 Ū	6 U
cis-1,3-Dichloropropene		6	U	6 U	6 L	1 6 U	6 U	6 Ū
Trichloroethene		6	U	6 U	6 t	1 6 U	140	121 %
Olbromochloromethane		6	U	6 U	6 L	6 U	6 U	6 Ü
		6	Ü	6 U	6 L	i 6 Ū	6 Ū	6 Ū
Benzene		6	U	6 U	6 t	i 6 U	6 Ū	95 %
Trans-1,3-Dichloroprope	ne	6	U	6 U	6 l	1 6 U	6 U	6 Ü
Bromoform		6	Ü	6 U	6 L	. 6 Ū	6 Ū	6 Ŭ
4-Methy1-2-pentanone		12	Ū	12 Ū	11 (1Ĭ Ū	11 Ŭ
2-Hexanone		12	Ū	12 U	11 (ii ū	ii ŭ
Tetrachloroethene			Ū	6 U	6 (• • •	
I C LI Q CII I UI UC LII CII C								
1,1,2,2-Tetrachloroetha	ne	. 6	U	6 Ū	6 i		6 Ü	6 U

RFW Batch Number: 9109G724		Client:	Ek	kco Houseware	_		W	ork Order: 29	94-	-02-03-0000		Page:	16
Cust	ID:	Trip Blank GW		R-1-GW		R-1	-GW	R-3-GW		R-3-GN		R-3-GW	
RFV	I# :	001		002		002	DL	003		003 MS		003 MSD)
Toluene		5	U	J 5	Ū	NA		5	Ü	87	%	86	*
Chlorobenzene		5	U	J 5	U	NA		5	U	91	%	90	%
Ethylbenzene		5	U	J 5	U	NA		5	U	5	Ü	5	Ü
Styrene		5	U	J 5	U	NA		5	U	5	Ū	5	Ū
Xylene (total)	-	5	U	J 5	U	NA		5	U	5	Ū	5	Ū
2-Chloroethylvinylether		10	U	J 10	U	NA		10	Ū	10	Ü	10	Ŭ
*= Outside of EPA CLP QC limits	.			56	9								
				,	•			108	,				

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Roy F. Weston, Inc. - Gulf Coast Laboratories

VOLATILES BY GC/MS, HSL LIST
Report Date: 10/10/91 13:07

RFW Batch Number: 9109G724		Client:	Ek	co H	<u>louseware</u>			Wor	<u>rk Order: 29</u>	94-	·02-03-0000		Page:	<u>la</u>
Cus	t ID:	Trip Blank			R-1-GW		R-1-G	W	R-3-6W		R-3-GW		R-3-GW	ı
Information Ma	RFW#: trix: D.F.: nits:	001 WATER	0		002 WATER 1.0 ug/L	0	002 D Water 2. ug/	00	003 WATER 1.0 ug/L	0	003 MS WATER 1.00 ug/L)	003 MSD WATER 1.0 ug/L	0
					-							- -		
Tolue Surrogate Bromofluorobe Recovery 1,2-Dichloroetha	nzen <mark>e</mark> ne-d4	102	% % %		99 100 107	% % % =f]	102 105 94	% % % ==f1	98 103 104	% % %	98 100 105	% % %	100 102 105	% % % :=f}
Chloromethane			U	===	10	- T I:	NA NA	T ;	10	-1 (10	- T I	10	U
			U		10	U	·NA		10	U	10	U	10	U
Bromomethane Vinyl Chloride		10	U		25		NA		10	U	10	U	10	U
Chloroethane		10	U		10	U	NA		10	U	10	U	10	U
Chloroethane Methylene Chloride		16			5	U	NA		5	U	26		25	
			U		10	U	NA		10	U	10	U	10	U
Acetone Carbon Disulfide		5	U		5	U	NA		5	U	5	U	5	U
1,1-Vichloroethene		5	U		5	U	NA		5	U	87	%	85	%
I.I-DICTIOTOETHANE		5	U		17		NA		51		48		47	
1,2-Dichloroethene (total)			U		83		NA		5	U	5	U	5	U
			U		5	U	NA		5	U	5	U	5	U
Chloroform 1,2-Dichloroethane			U		5	U	NA		5	U	5	U	5	U
			U		10	U	NA		10	U	10	U	10	U
Z-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride		5	U		54 -		NA		49		46		53	
Carbon Tetrachloride		- 5	U		5	U	NA		5	U	5	U	5	U
Vinyl Acetate		10	U		10	U	NA		10	U	10	U	10	U
Vinyl Acetate Bromodichloromethane		5	U		5	U	NA		5	U	5	U	5	U
			U		5	U	NA		5	U	5	U	5	U
1,2-Dichloropropane cis-1,3-Dichloropropene		5	U		5	U	NA		5	U	5	U	5	U
Trichloroethene		5	U			E	390 390		8		86	%	91	%
Dibromochloromethane	·	5	U		5	U	NA		5	U	5	U	5	U
1,1,2-Trichloroethane			U		5	U	NA		5	U	5	U	5	U
Benzene		5	U		5	U	NA		5	U	97	%	95	%
Trans-1,3-Dichloropropene		_ 5	U		5	U	NA		5	U	5	U	5	U
N		-	U		5	U	NA		5	U	5	U	. 5	Ū
dromoform_ 4-Methyl-2-pentanone		10	U		10	U	NA		10	U	10	U	10	U
Z-Hexanone		_ : 10	U		10	U	NA		10	U	10	U	10	U
Tetrachloroethene		5	U		5	U	NA		5	U	5	U	5	U
1,1,2,2-Tetrachloroethane		5	U		5	U	NA		5	U	5	U	5	ij
*= Outside of EPA CLP QC Tim	its.	-												

RFW Batch Number: 9109G745 Client: Ekco Houseware Work Order: 2994-02-03-0000 Page: 4b Cust ID: VBLK VBLK BS RFW#: 91GVE245-MB1 91GVE245-MB1 Toluene U % % 86 5 Chlorobenzene Ũ 90 5 U 5 U 5 U 10 U Ethylbenzene U Styrene

Xylene (total)
2-Chloroethylvinylether

*= Outside of EPA CLP QC limits. Š Ü 5 U 10 Ü

Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS, HSL LIST

Report Date: 10/15/91 14:05

RFW Batch Number: 9109G745	Client: Ek	co Houseware	
Cust ID:	VBLK	VBLK BS	· · · · · · · · · · · · · · · · · · ·
Sample RFW#:	91GVE245-MB1	91GVE245-MB1	
Information Matrix:	WATER	WATER	
D.F.:	1.00	1.00	
Units:	ug/L	ug/L	
toluene-d8			
Surrogate Bromofluorobenzene			
Recovery 1,2-Dichloroethane-d4			
Chloromethane			CERTERTS
Bromomethane	10 U	10 U	
Vinyl Chloride	10_U	10 U	
Lnioroethane	10 U		
Methylene Chloride	_ 5 U		
Acetone	_ 10 U		
Carbon Disulfide	5 U		
1,1-Dichloroethene	_ 5 U	7,0	
1,1-Dichloroethane	_ 5 U	• •	
1,2-Dichloroethene (total)	_ 5 U	• •	
Chloroform	5 U		
1,2-Dichloroethane	_ 5 U		
2-Butanone	10_U		
1,1,1-Trichloroethane	_ 5 U		
Carbon Tetrachloride		• •	
Vinyl Acetate Bromodichloromethane	- '! :		•
cis-1,3-Dichloropropene	- 5 U		
Trichloroethene	- 5 U		
Dibromochloromethane	_ 5 U		
1,1,2-Trichloroethane	- 5 U		
Benzene	- 5 U		
Trans-1,3-Dichloropropene	- 5 U		
Bromoform	- 5 Ü		
4-Methyl-2-pentanone	- 10 Ŭ	• •	
2-Hexanone	- : 10 Ū		
Tetrachloroethene	- ' 5 U		
1,1,2,2-Tetrachloroethane	5 U		
*= Outside of EPA CLP QC Timits.	-		

Cu	ist ID:	VBLK BS	O Houseware VBLK	VBLK BS	<u>k Order: 2994-</u> VBLK	VBLK BS	Page VBLK
	RFW#:	92GVE082-MB1	92GVE095-MB1	92GVE095-MB1			926VE092
Chlorobenzene		74 * %	0.1 U	85 %	0.1 U	205 * %	
1 1 1 2 Total Blancakland		_ ^ ^ 11	0.1 U	0.2 U	0.1 U		0.
Fthylhenzene		0.1 U	0.2 U	0.2 U	0.2 U		0.
Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane Isopropylbenzene	*	0.2 U	0.1 U	0.1 U	0.1 U 0.2 U	0.1 U	0.
p-Xvlene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U 0.2 U	0.
m-Xvlene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.
g-Xvlene		0.3 U	0.3 U	0.2 U	0.2 U	0.2 U	0.
Bromohenzene		0.1 U	0.5 U	0.3 U	0.3 U 0.1 U		0.
1.2.3-Trichloronronane		0.4 U	0.4 U	0.1 U	0.1 U	0.1 U 0.4 U	0. 0.
Isonronylhenzene		0.2 U	0.7 U	0.4 U	0.4 U		
n-Pronvilhenzene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U 0.2 U	0.
2-Chlorotoluene		0.2 U	0.2 U	0.2 U	0.2 U		0.
Isopropylbenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.
4-Chlorotoluene 1,3,5-Trimethylbenzene tert-Rutylbenzene		0.1 U	0.2 U			0.2 U	0.
tert-Butylbenzene		0.1 U	0.1 U	0.1 U 0.2 U	0.1 U 0.2 U	0.1 U	0.
1 2 4 Tulmakhulhamana			0.2 U	0.2 U		0.2 U	0.
sec-Rutylhenzene		0.3 U	0.3 U			0.3 U	0.
sec-Butylbenzene p-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene n-Butylbenzene		0.2 U	0.2 U	715	0.2 U	0.2 U	0.
1 3-Dichlersherzene		0.2 U	0.2 U 0.1 U	0.2 U 0.1 U	0.2 U	0.2 U	0.
1 4-Dichlerchenzene		0.1 U	0.1 U		0.1 U	0.1 U	0.
1 2-Dichlerohenzene		0.3 U	0.3 U 0.1 U		0.3 U	0.3 U	0.
n-But vi honzono		0.1 U	0.1 U		0.1 U	0.1 U	Ō.
n-Butylbenzene 1,2-Dibromo-3-Chloropropane 1,2,4-Trichlorobenzene Hexachlorobutadiene		0.1 U	0.1 U	0.1 U 0.5 U	0.1 U	0.1 U	0.
1,2-010f0m0-3-chiofopfopane		0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.
Hovachlarabutadiana	·	0.2 U	0.4 U	0.2 U	0.2 U	0.2 U	0.
Nanhthalana		0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.
1 2 2 Trichlarchenzana		0.2 U	0.2 U		0.2 U	0.2 U	0.
cie 1 3 Dichlamonyonen		0.1 U			0.1 U	0.1 U	0.
Naphthalene 1,2,3-Trichlorobenzene cis-1,3-Dichloropropene trans-1,3-Dichloropropene		0.1 U		0.1 U	0.1 U	0.1 U	0.
Acetone		0.3 U 2 U		0.3 U	0.3 U	0.3 U	0.
Acetone Carbon Disulfide		0.1 U	2 U 0.1 U	2 U	2 U	2 U	•
2-Butanone		0.1 U	0.1 U 0.3 U	0.1 U	0.1 U	0.1 U	0.
2-Butanone Vinyl Acetate 4-Methyl-2-pentanone		0.3 U	0.3 U	0.3 U 0.3 U	0.3 U	0.3 U	0.
A. Mathyl - 2 - nontanona		0.3 U	• • • •		0.3 U	0.3 U	0.
7 Hermone	· ·	0.4 U	• • • -		0.4 U	0.4 U	0.
2-Hexanone 2-Chloroethylvinylether		0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	. 0.
*= Outside of EPA CLP QC 1i	=11.	. 0.2 0	0.2 U	0.2 U	0.2 U	0.2 U	0.

Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS, METHOD 524 LIST
Client: Ekco Houseware Work Order: 24 LIST Report Date: 04/15/92 11:21 Work Order: 2994-02-03-0000 Page: 7a RFW Batch Number: 9203G478 Page: 7a

	Cust ID:	VBLK BS	VBLK	VBLK BS	VBLK	VBLK BS	VBLK (
Sample		92GVE082-MB1	92GVE095-MB1	92GVE095-MB1	92GVE083-MB1	92GVE083-MB1	92GVE092-MB1
Information	Matrix:	WATER	WATER	WATER	WATER	WATER	WATER
	D.F.:	1.00	1.00	1.00	1.00	1.00	1.00
	Units:	UG/L	UG/L	UG/L	UG/L	UG/L	UG/L
Bromofluo	robenzene	110 %	111 %	111 %	100 %	93 %	111 %
Dichlorodifluoromethane			0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Chloromethane			0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Bromomethane		กาแ	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Vinyi Unioride		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Chloroethane		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Chloroethane Methylene Chloride Trichlorofluoromethane		0.6 U	0.6 U	0.6 U	0.6 U	0.6 U	0.6 U
Trichlorofluoromethane		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,1-Dichloroethene		92 %	0.1 U	113 %	0.1 U	259 * %	0.1 U
1,1-Dichloroethane		0.1 U	0.1 U	0.1 U	0.1 U	0.1 Ü	0.1 U
1,1-Dichloroethane cis-1,2-Dichloroethene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
2,2-Dichloropropane	_	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
2,2-Dichloropropane trans-1,2-Dichloroethene		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Chloroform		0211	0.2 U	0.2 U	0.2 U	0.2 U	0.2 Ü
Bromochioromethane		0211	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,2-Dichloroethane 1,1,1-Trichloroethane Carbon Tetrachloride		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,1,1-Trichloroethane		0.2 U	0.2 · U	0.2 U	0.2 U	0.2 U	0.2 U
Carbon Tetrachloride		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,1-Dichloropropene		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,1-Dichloropropene		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Dibromomethane		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 Ū
Dibromomethane 1,2-Dichloropropane		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Trichloroethene		120 %	0.3 U	136 * %	0.3 U	266 * %	0.3 U
Trichloroethene Dibromochloromethane		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,2-Dibromoethane 1,1,2-Trichloroethane		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,1,2-Trichloroethane		0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Delizelle	•	111 74	0.2 U	81 %	0.2 U	172 * %	0.2 U
1,3-Dichloropropane	·	0.2 Û	0.2 U	0.2 Ü	0.2 U	0.2 Ü	0.2 U
Bromoform		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Tetrachloroethene		0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,1,2,2-Tetrachloroethan	e		0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Toluene	-	69 * %	0.1 U	76 %	0.1 U	171 * %	0.1 U
*= Outside of EPA CLP QC	limits.	_	 0	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	0.1 0	211 /6	0.1 0

RFW Batch Number: 9203G478 Cust ID:	Client: Ekc S-12-GW		Wo	<u>rk Order: 2994-</u>	02-03-0000	Page: 6b
cust iv:	3-12-GW	S-12-GW	5-12-GW	VBLK	VBLK BS	VBLK
RFW#:	020 DL	020 MS	020 MSD	92GVE081-MB1	92GVE081-MB1	92GVE082-MB1
Chlorobenzene 1,1,1,2-Tetrachloroethane	NA	198 * %	219 * %	0.1 U	156 * %	0.1 U
1,1,1,2-Tetrachloroethane	NA	0.2 U	0.2 U	0.2 U	0.2 Ü	0.2 U
Ethylbenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Tilizate rachioroechane Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 Ū
p-Xylene	NA	0.2 U	0.2 U		0.2 U	0.2 U
m-Xylene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
p-Xylene	·NA	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Bromobenzene	NA	0.1 U	0.1 U		0.1 U	0.1 U
1,2,3-Trichloropropane	NA	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
Isopropylbenzene n-Propylbenzene 2-Chlorotoluene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
n-Propylbenzene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
?-Chlorotoluene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
I-Chlorotoluene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,3,5-Trimethylbenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
tert-Butylbenzene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
tert-Butylbenzene	NA	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
sec-Butylbenzene b-Isopropyltoluene l,3-Dichlorobenzene l,4-Dichlorobenzene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
o-Isopropyitoluene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,3-Dichlorobenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
.4-Dichlorobenzene	NA	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
1.2-Dichlorobenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
n-Butylbenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
n-Butylbenzene 1,2-Dibromo-3-Chloropropane 1,2,4-Trichlorobenzene	NA	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
.2.4-Trichlorobenzene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
		0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
laphthalene 1,2,3-Trichlorobenzene :is-1,3-Dichloropropene :rans-1,3-Dichloropropene	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
.2.3-Trichlorobenzene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
is-1.3-Dichloropropene	NA	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
rans-1.3-Dichloropropene	NA	0.3 U	0.3 U	0.3 U	0.3 U	0.3 Ü
cetone	NA	2 U	2 U	2 U	2 U	2 U
cetone arbon Disulfide	NA	0.1 U	0.1 Ŭ	0.1 Ū	0.1 U	0.1 Ŭ
-Butanone	NA	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
invl Acetate	NA	0.3 U	0.3 U		0.3 U	0.3 U
P-Butanone Vinyl Acetate I-Methyl-2-pentanone	NA	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
/-Heyanone	NA	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
2-Chloroethylvinylether	NA	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
- Outside of EPA CLP QC limits.	• • •					•

18-78

RFW Batch Number:	9203 G 478	VOLAT Client: Ekco	ILES BY GC/MS, Houseware	METHOD 524 I	LIST <mark>rk Order: 2994-</mark>	Report Date: 02-03-0000	04/15/92 11:21 Page: 6a
	Cust ID:	S-12-GW	S-12-GW	S-12-GW	VBLK	VBLK BS	VBLK (
Sample	RFW#:	020 DL	020 MS	020 MSD	92GVE081-MB1		926VE082-MB1
Information	Matrix:	WATER	WATER	WATER	WATER	WATER	WATER
	D.F.: Units:	100 UG/L	1.00 UG/L	1.00 UG/L	1.00 UG/L	1.00 UG/L	1.00 UG/L
Bro	mofluorobenzene	112 %	93 %	95 %		100 %	
Dichlorodifluorome	======================================	f1- NA	f1-: 0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Chloromethane	LII GII C	NA NA	0.3 U	0.3 U		0.3 U	0.3 U
Chloromethane Bromomethane		NA NA	0.1 U	0.1 U		0.1 U	0.1 U
Bromomethane Vinyl Chloride		NA NA	5	0.2 U	0.2 U	0.2 U	0.2 U
Chloroethane		NA NA	0. Ĭ U	0.1 U	0.1 U	0.1 U	0.1 U
Chloroethane Methylene Chloride Trichlorofluoromet	·····	NA NA	8	8	0.6 U	0.6 U	0.6 U
Trichlorofluoromet	hane	NA	0.Ĭ U	0.Ĭ U	0.1 U	0.1 U	0.1 U
1,1-Dichloroethene		NA	279 * %	268 * %	0.1 U	201 * %	0.1 U
1,1-Dichloroethane		NA	0.1 Ü	0.1 U	0.1 U	0.1 U	0.1 U
1,1-Dichloroethane cis-1,2-Dichloroet	hene	NA	E	70	0.2 U	0.2 U	0.2 U
2,2-Dichloropropand	е	NA	0.1 Ū	0.1 U	0.1 U	0.1 U	0.1 U
trans-1.2-Dichloro	ethene	NA	0.1 U	0.1 U		0.1 U	0.1 U
Chloroform		NA	0.2 U	0.2 U		0.2 U	0.2 U
Chloroform Bromochloromethane 1,2-Dichloroethane 1,1,1-Trichloroeth Carbon Tetrachlori		NA	0.2 U	0.2 U	• • • •	0.2 U	
1,2-Dichloroethane		NA	0.2 U	0.2 U	7.2	0.2 U	
1,1,1-Trichloroeth	ane	NA	5	4	0.2 U	0.2 U	
Carbon Tetrachloric	de	NA	0.1 U	0.1 U		0.1 U	
1,1-Dichloropropen	e	NA	0.1 U	0.1 U	* • • •	0.1 U	4.5
1,1-Dichloropropend Bromodichlorometha	ne	NA	0.2 U	0.2 U		0.2 U	7.5
vibromomethane		NA NA	0.2 U	0.2 U	·	0.2 U	• · • -
1,2-Dichloropropan	e	NA	0.2 U	0.2 U		0.2 U	
Trichloroethene		2300	732 * %	0 * %	0.3 U	219 * %	0.3 U

0.2 U

0.2 U

0.3 U

177 * %

0.2 U

0.2 U

0.1 U

0.2 U

170 * %

NA

NA

NA

NA

NA

NA

NA

NA

NA

0.2 U

0.2 U

0.3 U

0.2 U

0.2 U

0.2 U

0.1 U

0.2 U

0.1 U

0.2 U

0.2 U

0.3 U

0.2 U

0.2 U

0.1 U 0.2 U

179 * %

0.2 U

0.2 U

0.3 U

157 * %

0.2 U

0.2 U

0.1 U

0.2 U

132 * %

0.2 U

0.2 U

0.3 U

0.2 U

0.2 U

0.2 U

0.1 U

0.2 U

0.1 U

Dibromochloromethane

1,1,2-Trichloroethane

1,1,2,2-Tetrachloroethane

*- Outside of EPA CLP QC limits.

1,3-Dichloropropane

Tetrachloroethene

1,2-Dibromoethane

Benzene

Toluene

Bromoform

RFW Batch Number: 9203G478	<u> Client: Ekc</u>		<u>Work</u>	Order: 2994-0		Page:
Cust ID:	R-12-GW	R-12-GW	S-4-GW	S-11-GW	S-12-GW	S-12-GW
RFW#:	017	017 DL	018	019	020	020 DL
Chlorobenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
Chlorobenzene 1,1,2-Tetrachloroethane	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
Ethylbenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA NA
Styrene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA NA
p-Xylene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
m-Xylene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
o-Xylene	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
Bromobenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
1,2,3-Trichloropropane	0.4 U	NA	0.4 U	0.4 U	0.4 U	NA
Isopropylbenzene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
n-Propylbenzene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
2-Chlorotoluene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
4-Chlorotoluene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
1,3,5-Trimethylbenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
tert-Butylbenzene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
1,2,4-Trimethylbenzene	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
sec-Butylbenzene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
p-Isopropyltoluene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
1,3-Dichlorobenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
1,4-Dichlorobenzene	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
1,2-Dichlorobenzene	0.1 U	NA	0.1 U	0.1 U	0.1 Ü	NA
n-Buty1benzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
1,2-Dibromo-3-Chloropropane	0.5 U	NA	0.5 U	0.5 U	0.5 Ü	NA
1,2,4-Trichlorobenzene	0.2 U	NA .	0.2 U	0.2 U	0.2 U	NA
Hexachlorobutadiene	0.4 U	NA	0.4 U	0.4 U	0.4 U	NA
Naphthalene	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
1,2,3-TrichTorobenzene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
cis-1,3-Dichloropropene	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
trans-1,3-Dichloropropene	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
Acetone	2 U	NA	2 U	2 U	2 U	NA
Carbon Disulfide	0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
Chlorobenzene 1,1,1,2-Tetrachloroethane Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane Isopropylbenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene 1,3,5-Trimethylbenzene tert-Butylbenzene 1,2,4-Trimethylbenzene sec-Butylbenzene p-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene 1,2-Oibromo-3-Chloropropane 1,2,4-Trichlorobenzene Hexachlorobutadiene Naphthalene 1,2,3-Trichlorobenzene cis-1,3-Dichloropropene trans-1,3-Dichloropropene Acetone Carbon Disulfide 2-Butanone Vinyl Acetate 4-Methyl-2-pentanone 2-Hexanone 2-Chloroethylvinylether	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
Vinyl Acetate	0.3 U	NA	0.3 U	0.3 U	0.3 U	NA
4-Methy1-2-pentanone	0.4 U	NA	0.4 U	0.4 U	0.4 U	NA
2-Hexanone	0.5 U	NA	0.5 U	0.5 U	0.5 U	NA
2-Chloroethylvinylether	0.2 U	NA	0.2 U	0.2 U	0.2 U	NA

Cust ID: R-12-GW R-12-GW S-4-GW S-11-GW S-12-GW S	RFW Batch Number: 9203G47	78	VOLAT Client: Ekco	ILES BY GC/MS, Houseware	METHOD 524 LIS Work	ST Order: 2994-0	Report Date: 0 ⁴ 2-03-0000	4/15/92 11:21 Page: 5a
Sample		Cust 1D:	R-12-GW	R-12-GW	S-4-GW	S-11-GW	S-12-GW	
D.F.: 1.00 2.00 1.00 1.00 1.00 10.0						019	020	-
D.F.: 1.00 2.00 1.00	Information		WATER	WATER	WATER	WATER	WATER	WATER
Bromofluorobenzene			1.00	2.00	1.00	1.00	1.00	10.0
Dichlorodifluoromethane		Units:	UG/L	UG/L	UG/L	UG/L	UG/L	UG/L
Dichlorodifiluoromethane					96 %		92 %	115 %
Chloromethane	Dichlorodifluoromethane		0.3 U				• •	• •
Bromomethane	Chloromethane		0.2 U	NA	0.2 U			
Viny1 Chloride	Bromomethane							
Chloroethane	Vinyl Chloride		0.2 U		- · · - ·		-	
	Chloroethane		0.1 U			_	0.1 U	
	Methylene Chloride		9	NA	" u	8	9	NA
	Trichlorofluoromethane		2	NA		2		
1,1-Dichloroethane	I,1-Dichloroethene		3		: I I		0.1 U	
cis-1,2-Dichloroethene 16 NA 0.2 U 0.2 U E 36 2,2-Dichloropropane 0.1 U NA 0.1 U 0.1 U 0.1 U 0.1 U NA trans-1,2-Dichloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA Bromochloromethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dichloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,1,1-Trichloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,1-Dichloropethane 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1-Dichloropropene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1-Dichloropropene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dibromoethane 0.2 U NA 0.2 U </td <td>1,1-Dichloroethane</td> <td></td> <td>E</td> <td>22</td> <td>0.1 U</td> <td></td> <td></td> <td></td>	1,1-Dichloroethane		E	22	0.1 U			
2,2-Dichloropropane	cis-1,2-Dichloroethene	· · · · · · · · · · · · · · · · · · ·	16	NA	0.2 U		E	36
trans-1,2-Dichloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA Chloroform 0.2 U NA 0.2 U 0.2 U 0.2 U NA Bromochloromethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dichloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,1,1-Trichloroethane 0.1 U NA 0.1 U 0.1 U 0.1 U NA Carbon Tetrachloride 0.1 U NA 0.1 U 0.1 U 0.1 U NA Bromodichloropropene 0.1 U NA 0.1 U 0.1 U 0.1 U NA Dibromomethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dibromoethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dibromoethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,2-Dibromoethane 0.2 U NA 0.2 U 0.2 U <t< td=""><td>2,2-Dichloropropane</td><td></td><td>0.1 U</td><td>NA</td><td>0.1 U</td><td>0.1 U</td><td>0.1 Ü</td><td>NA</td></t<>	2,2-Dichloropropane		0.1 U	NA	0.1 U	0.1 U	0.1 Ü	NA
Chloroform	trans-1,2-Dichloroethene	-	0.1 U	NA	0.1 U		0.1 U	
Bromochloromethane	Chioroform		N 2 II	NA	0.2 U	0.2 U	0.2 U	
	Bromochloromethane		0.2 U .	NA	0.2 U	0.2 U		
	1,2-Dichloroethane		0.2 U			0.2 U	0.2 U	
	1,1,1-Trichloroethane		0.2 U	NA .	0.2 U	0.2 U	. 5	NA
	Carbon Tetrachloride		0.1 U	NA	0.1 U	0.1 U	0.1 U	NA
1,2-Dichloropropane	1.1-Dichloropropene		0.1 U	NA	0.1 U	0.1 U		
1,2-Dichloropropane	Bromodichloromethane		0.2 U	NA	0.2 U	0.2 U	0.2 U	NA
1,2-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA Trichloroethene 17 NA 5 ; 0.3 U E E Dibromochloromethane 0.2 U NA 0.2 U 0.2 U 0.2 U 0.2 U NA 1,2-Dibromoethane 0.2 U NA 0.2 U 0.2 U 0.2 U 0.2 U NA 1,1,2-Trichloroethane 0.3 U NA 0.3 U 0.3 U 0.3 U 0.3 U NA Benzene 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,3-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA Bromoform 0.2 U NA 0.2 U 0.2 U 0.2 U NA Tetrachloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1,2,2-Tetrachloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U 0.2 U NA	Dibromomethane		0.2 U	NA	0.2 U	0.2 U		NA
Trichloroethene	1.2-Dichloropropane		0.2 U					
1,1,2-Trichloroethane 0.2 U NA 0.2 U 0.2 U 0.3 U NA 1,1,2-Trichloroethane 0.3 U NA 0.3 U 0.3 U 0.3 U NA Benzene 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,3-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA Bromoform 0.2 U NA 0.2 U 0.2 U 0.2 U NA Tetrachloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1,2,2-Tetrachloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA	Trichloroethene		17			0.3 U	E	E
1,2-Dibromoethane 0.2 U NA 0.2 U 0.2 U 0.3 U NA 1,1,2-Trichloroethane 0.3 U NA 0.3 U 0.3 U 0.3 U NA Benzene 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,3-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA Bromoform 0.2 U NA 0.2 U 0.2 U 0.2 U NA Tetrachloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1,2,2-Tetrachloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA	Dibromochloromethane		0.2 U				0.2 Ũ	NA
1,1,2-Trichloroethane 0.3 U NA 0.3 U 0.3 U 0.3 U NA Benzene 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,3-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U 0.2 U NA Bromoform 0.2 U NA 0.2 U 0.2 U 0.2 U 0.2 U NA Tetrachloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1,2,2-Tetrachloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA	1.2-Dibromoethane		0.2 U	NA	0.2 U			NA
Benzene 0.2 U NA 0.2 U 0.2 U 0.2 U NA 1,3-Dichloropropane 0.2 U NA 0.2 U 0.2 U 0.2 U NA Bromoform 0.2 U NA 0.2 U 0.2 U 0.2 U NA Tetrachloroethene 0.1 U NA 0.1 U 0.1 U 0.1 U NA 1,1,2,2-Tetrachloroethane 0.2 U NA 0.2 U 0.2 U 0.2 U NA	1.1.2-Trichloroethane		0.3 U		0.3 U	0.3 U	0.3 U	NA
### ##################################	Benzene		0.2 U					
### ##################################	1,3-Dichloropropane		0.2 U					NA
l,l,2,2-Tetrachloroethane O.2 U NA O.2 U O.2 U O.2 U NA	Bromotorm		0.2 U		0.2 U		0.2 U	NA
l,l,2,2-Tetrachloroethane O.2 U NA O.2 U O.2 U O.2 U NA	Tetrachloroethene			NA	0.1 U	0.1 U	0.1 U	NA
Toluene O.1 U NA O.1 U O.1 U NA	1,1,2,2-Tetrachloroethane	9						
	Toluene		0.1 U					

RFW Batch Number: 9203G478	Client: Ekco		Work	Order: 2994-		Page: 4b
Cust ID:	I-4-GW	I -4-GW	R-7-GW	R-7-GW	R-10-GW	R-10-GW
RFW#:	014	014 DL	015	015 DL	016	016 DL 🚚
Chlorobenzene	0.1 U	NA NA	0.1 U	NA	0.1 U	NA C
1,1,1,2-Tetrachloroethane		NA	0.2 U	NA	0.2 U	NA T
I,I,I,Z-Tetrachioroethane Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane Isopropylbenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
Styrene	0.2 U	NA	0.2 U	NA	0.2 U	NA
p-Xylene	0.2 U	NA	0.2 U	NA	0.2 U	NA
m-Xylene	0.2 U	NA	0.2 U	NA	0.2 U	NA
o-Xylene	0.3 U	NA	0.3 U	NA	0.3 U	NA
Bromobenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
1,2,3-Trichloropropane	0.4 U	NA	0.4 U	NA	0.4 U	NA
1,2,3-1richioropropane Isopropylbenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene 1,3,5-Trimethylbenzene tert-Butylbenzene 1,2,4-Trimethylbenzene sec-Butylbenzene	0.2 U	NA	0.2 U	NA	0.2 U	NA
n-Propylbenzene	0.2 U	NA	0.2 U	NA	0.2 U	NA
2-Chlorotoluene	0.2 U	NA	0.2 U	NA	0.2 U	NA
4-Chlorotoluene	0.2 U	NA	0.2 U	NA	0.2 U	NA
1,3,5-Trimethylbenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
tert-Butylbenzene	0.2 U	NA	0.2 U	NA	0.2 U	NA
1,2,4-Trimethylbenzene	0.3 U	NA	0.3 U	NA	0.3 U	NA
sec-Butylbenzene p-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene	0.2 U	NA	0.2 U	NA	0.2 U	NA
p-Isopropyltoluene	0.2 U	NA	0.2 U	NA	0.2 U	NA .
1,3-Dichlorobenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
1,4-Dichlorobenzene	0.3 U	NA	0.3 U	NA	0.3 U	NA
1,2-Dichlorobenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
n-Butylbenzene	0.1 U	NA	0.1 U	NA	0.1 U	NA
1,2-Dibromo-3-Chloropropane	0.5 U	NA	0.5 U	NA	0.5 U	NA
n-Butylbenzene 1,2-Dibromo-3-Chloropropane 1,2,4-Trichlorobenzene Hexachlorobutadiene	0.2 U	NA .	0.2 U	NA	0.2 U	NA
Hexachlorobutadiene	0.4 U	NA	0.4 U	NA	0.4 U	NA
Naphthalene	0.2 U	NA	0.2 U	NA	0.2 U	NA
Naphthalene 1,2,3-Trichlorobenzene cis-1,3-Dichloropropene trans-1,3-Dichloropropene	0.1 U	NA	0.1 U	NA	0.1 U	NA
cis-1,3-Dichloropropene	0.1 U	NA	0.1 U	NA	0.1 U	NA
trans-1,3-Dichloropropene	0.3 U	NA	0.3 U	NA	0.3 U	NA
Acetone	2 U	NA	2 U	NA	2 U	NA
Acetone Carbon Disulfide	0.1 U	NA	0.1 U	NA	0.1 V	NA
2-Butanone	0.3 U	NA	0.3 U	NA	0.3 U	NA
Vinyl Acetate	0.3 U	NA	0.3 U	NA	0.3 U	NA
2-Butanone Vinyl Acetate 4-Methyl-2-pentanone	0.4 U	NA	0.4 U	NA	0.4 U	NA
2-Hexanone	0.5 U	NA	0.5 U	NA	0.5 U	NA
2-Chloroethylvinylether	0.2 U	NA	0.2 U	NA	0.2 U	NA
*- Outside of EPA CLP QC limits.						

RFW Batch Number: 9203G478		V Client:	OLAT Ekco	ILES BY GC/MS Houseware	, METHOD 5		ST Order: 2994-0	Report Date: 0 2-03-0000	4/15/92 11:21 Page: 4a
	ust ID:	1-4-6W		1-4-GW	R-7-G		R-7-GW	R-10-6W	R-10-6W
		1-4-4		1-4-GW	K-7-0		K-/-G#	K-10-0#	4-10-0M (.
Sample	RFW#:	014		014 DL	019		015 DL	016	016 DL
Information)	latrix:	WATER		WATER	WATER		WATER	WATER	WATER
	D.F.:	1.0	0	25.0	1.0		5.00	1.00	10.0
	Units:	UG/L		UG/L	UG/1	L	UG/L	UG/L	UG/L
Bromofluorob	enzene	106	%	117 %	99	%	113 %	94 %	115 %
Dichlorodifluoromethane		0.3	-71 U	f1 Na	0.3	f1-: U	f1- NA	fl- 0.3 U	f1 NA
Chloromethane		0.2	Ŭ	NA NA	0.3		NA NA	0.3 U 0.2 U	NA NA
			Ŭ	NA NA	0.1	ŭ	NA NA	0.2 U	NA NA
Bromomethane Vinyl Chloride		0.2	Ŭ	NA NA	0.2	Ü	NA	35	NA NA
Chloroethane		0.1	Ŭ	NA NA	0.1	Ŭ	NA NA	0.1 U	NA NA
Chloroethane Methylene Chloride Trichlorofluoromethane		8	u	NA	7	•	NA	8	NA NA
Trichlorofluoromethane		0.1	Ŭ`	NA NA	0.1	U	NA	0.1 U	NA NA
1,1-Dichloroethene		0.1	Ŭ	NA	7		NA	27	NA NA
1.1-Dichloroethane			Ě	K96	· ·	E	38	F	260
cis-1,2-Dichloroethene		0.2	Ū	NÁ	6	_	NA	Ē	110
			Ū	NA	0.1	U	NA	0.1 Ū	NA
2,2-Dichloropropane trans-1,2-Dichloroethene		0.1	Ū	NA	0.1	Ŭ	NA	0.1 U	NA
Chloroform		0.2	Ū	NA	0.2	Ŭ	NA	0.2 U	NA
Bromochloromethane		0.2	U	NA	0.2	U	NA	0.2 U	NA
1.2-Dichloroethane		ก <i>ว</i>	U	NA	0.2	U	NA	0.2 U	NA
1,1,1-Trichloroethane		0.2	U	NA .	6		NA	4	NA
Larbon letrachioride		11 1	U	NA	0.1	U	NA	0.1 U	NA
1,1-DichloropropeneBromodichloromethane		0.1	U	NA	0.1	U	NA	0.1 U	NA
Bromodichloromethane		0.2	U	NA	0.2	U	NA	0.2 U	NA
HINCOMOMOTRANO		11 2	U	NA	0.2	U	NA	0.2 U	NA
1,2-Dichloropropane		0.2	U	NA	0.2	U	NA	0.2 U	NA
I TI CRIATAGI BARA		T-7.7	•	NA		E	69	E	170
Dibromochloromethane		0.2	U	NA	0.2	U	NA	0.2 U	NA
1,2-Dibromoethane		0.2	U	NA	0.2	U	NA	0.2 U	NA
l,2-Dibromoethane l,1,2-Trichloroethane		0.3	U	NA	0.3	_	NA	0.3 U	NA
Benzene		0.2	U	NA	0.2	U	NA	0.2 U	NA
1,3-Dichloropropane		0.2	U	NA	0.2	_	NA	0.2 U	NA
Bromoform		0.2	U	ŅA	0.2	U	NA	0.2 U	NA
Tetrach1oroethene		0.1	U	NA	0.1	U	NA	0.1 U	NA
1,1,2,2-Tetrachloroethane		0.2	U	NA	0.2		NA	0.2 U	NA
Toluene		0.1	IJ	NA	0.1	IJ	NA	0.1 U	NA
*- Outside of EPA CLP QC 11	mits.								

RFW Batch Number: 9203G47	Cust ID:	S-12-GW Dup	co <u>Houseware</u> S-12-GW Dup	S-12-FB	rk Order: 2994-02 S-12-Pump EB	P-3-GW	Page: 3b
			o iz an bap	3 12 15	3-12-1 dilly ED	r-3-un	r-3-aw
	RFW#:	010 DL	010 DL	011	012	013	013 DL
hlorobenzene ,1,1,2-Tetrachloroethane		NA	NA NA	0.1 U	ō.1 U	0.1 U	NA NA
,1,1,2-Tetrachloroethane		- NA	NA	0.2 U	0.2 U	0.2 U	NA
,1,1,2-letrachioroethane thylbenzene_ tyreneXyleneXylene -Xylene romobenzene_ ,2,3-Trichloropropane_ sopropylbenzene		NA	NA	0.1 U	0.1 U	0.1 U	NA
tyrene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
-Xylene_		- NA	NA	0.2 U	0.2 U	0.2 U	NA
-Xylene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
-Xylene		- NA	NA	0.3 U	0.3 U	0.3 U	NA
romobenzene		- NA	NA	0.1 U	0.1 U	0.1 Ŭ	NA
,2,3-Trichloropropane		NA NA	NA	0.4 U	0.4 U	0.4 U	NA NA
sopropylbenzene		- NA	NA	0.2 U	0.2 U	0.2 Ü	NA
sopropylbenzene -Propylbenzene -Chlorotoluene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
			NA	0.2 U	0.2 U	0.2 U	NA
·Chlorotoluene		NA NA	NA	0.2 U	0.2 U	0.2 U	NA
-Chlorotoluene ,3,5-Trimethylbenzene		- NA	NA	0.1 U	0.1 U	0.1 U	NA
ert-Butylbenzene		NA NA	NA	0.2 U	0.2 U	0.2 U	NA
2.4-Trimethylbenzene		- NA	NA	0.3 U	0.3 U	0.3 U	NA
ec-Butylbenzene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
Isopropyltoluene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
3-Dichlorobenzene		- NA	NA	0.1 U	0.1 U	0.1 U	NA
4-Dichlorobenzene		- NA	NA	0.3 U	0.3 U	0.3 U	NA
4-Dichlorobenzene		NA NA	NA	0.1 U	0.1 U	0.1 U	NA
Butylbenzene		- NA	NA	0.1 U	0.1 U	0.1 U	NA
-Butylbenzene ,2-Dibromo-3-Chloropropa ,2,4-Trichlorobenzene	ne	NA	NA	0.5 U	0.5 U	0.5 U	NA
2.4-Trichlorobenzene		- NA	NA .	0.2 U	0.2 U	0.2 U	NA
exacniorodutadiene		NA	* NA	0.4 U	0.4 U	0.4 U	NA
phthalene 2,3-Trichlorobenzene		- NA	NA	0.2 U	0.2 U	0.2 U	NA
2.3-Trichlorobenzene		- NA	NA	0.1 U	0.1 U	0.1 U	NA
s-1.3-Dichloropropene		- NA	NA	0.1 U	0.1 Ū	0.1 U	NA
s-1,3-Dichloropropene rans-1,3-Dichloropropene		- NA	NA	0.3 U	0.3 U	0.3 U	NA
etone	·	- NA	NA	2 U	2 Ü	2 U	NA
etone urbon Disulfide		- NA	NA NA	0.Î Ü	0. Ī Ŭ	0.1 Ŭ	NA
Butanone	·	- NA	NA	0.3 U	0.3 U	0.3 U	NA NA
nvl Acetate		- NA	NA NA	0.3 U	0.3 U	0.3 U	NA NA
Butanone Inyl Acetate Methyl-2-pentanone		- NA	NA NA	0.4 U	0.4 U	0.4 U	NA NA
-Hexanone	<u> </u>	- NA	NA NA	0.5 U	0.5 U	0.5 U	NA NA
Hexanone Chloroethylvinylether		- NA	NA NA	0.2 U	0.3 U	0.2 U	NA NA
- Outside of EPA CLP QC	limite	_ '''	1 1/1	V.E U	J.E U	V.E V	11/1

DEU Datah Number 200004	170	VOLA	ton, Inc Gul	METHOD 524 I	_IST R	eport Date: 04	
RFW Batch Number: 9203G4	1/8	Client: EKC	o Houseware	WOI	<u>rk Order: 2994-02</u>	-03-0000	Page: 3a
	Cust ID:	S-12-GW Dup	S-12-GW Dup	S-12-FB	S-12-Pump EB	P-3-GW	P-3-GW
Sample	RFW#:	010 DL	010 DL	011	012	013	013 DL
Information	Matrix:	WATER	WATER	WATER	WATER	WATER	WATER
	D.F.:	2.00	500	1.00	1.00	1.00	5.00
	Units:	UG/L	UG/L	UG/L	UG/L	UG/L	UG/L
	robenzene	107 %	108 %	107 %	109 %	99 %	111 %
Dichlorodifluoromethane_		NA NA	NA	0.3 U	0.3 U	f1 0.3 U	f1 NA
			NA	0.2 U	0.2 U	0.2 U	NA NA
———		-	NA NA	0.1 U	0.1 U	0.1 U	NA NA
Vinyl Chloride		- NA	NA	0.2 U	0.2 U	0.2 U	NA
Chloroethane	 	- NA	NA NA	0.1 U	0.2 U	0.2 U	NA NA
Chloroethane Methylene Chloride Trichlorofluoromethane		- NA	NA NA	4	3	0.6 U	NA NA
Irichlorofluoromethane		- NA	NA NA	0.7	0.5	0.0 U	NA NA
1.1-Dichloroethene		- NA	NA NA	0.1 U	0.1 U	6	NA NA
1,1-Dichloroethane		- NA	NA NA	0.1 U	0.1 U	E	84
cis-1,2-Dichloroethene_			NA NA	0.2 U	0.2 U	6	NA
2,2-Dichloropropane	 	- NA	NA NA	0.1 U	0.1 U	0.Ĭ U	NA NA
trans-1,2-Dichloroethene		- NA	NA	0.1 U	Ŏ. 1 Ŭ	0.1 Ŭ	NA
			NA	0.2 U	0.2 U	0.2 Ū	NA
			NA	0.2 U	0.2 U	0.2 U	NA
		•	NA	0.2 U	0.2 U	0.2 U	NA
1.1.1-Trichloroethane		- NA	NA ·	0.2 U	0.2 U	F	92
Carbon Tetrachloride		- NA	NA	0.1 U	0.1 U	0.1 Ū	NA
I.I-IJICDIOTODTODPOP		NΔ	NA	0.1 U	0.1 U	0.1 U	NA
Bromodichloromethane	"	- NA	NA	0.2 U	0.2 U	0.2 U	NA
Dibromomethane		- NA	NA	0.2 U	0.2 U	0.2 U	NA
1,2-Dichloropropane			NA	0.2 U	0.2 U	0.2 U	NA NA
Trichloroethene		- '''' E	2000	0.8	0.3 U	F	62
Dibromochloromethane		- NA	NA	0.2 U	0.2 U	0.2 บั	NA
1 2-Dibromoethane		- NA	NA NA	0.2 U	0.2 U	0.2 U	NA NA
1 1 2 Trichloroethane		- NA	NA	0.3 U	0.2 U	0.3 U	NA NA
1,2-Dibromoethane 1,1,2-Trichloroethane Benzene		- NA	NA NA	0.3 U	0.3 U 0.2 U	0.3 U	NA NA
1,3-Dichloropropane			NA NA	0.2 U	0.2 U	0.2 U	NA NA
Bromoform		- NA	NA NA	0.2 U	0.2 U	0.2 U	NA NA
Tetrachloroethene		- NA	NA NA	0.1 U	0.2 U	0.1 U	NA NA
1,1,2,2-Tetrachloroetham	<u> </u>	- NA	NA	0.2 U	0.1 U	0.2 U	NA NA
1,1,2,2-letrachionoethan Toluene		- NA	NA NA	0.2 U	0.2 U	0.2 U	NA NA
*= Outside of EPA CLP QC	11=11=	_ 11/4	IVA	0.1 0	0.1 0	0.1 0	IIA

RFW Batch Number: 9203G478	Client: Ekco		Wo	rk Order: 2994-		Page: 2b
Cust 1D:	I-13-GW	I-13-GW	I-14-GW	Trip Blank A	Trip Blank B	S-12-GW Dup
RF₩#:	006 MS	006 MSD	007	008	009	010
Chlorobenzene	162 * %	160 * %	0.1 U	0.1 U	0.1 U	0.1 U
1,1,1,2-Tetrachloroethane		0.2 Ü	0.2 U		0.2 U	0.2 U
Ethylbenzene	0.1 U	0.1 U	0.1 U		0.1 U	0.1 U
Styrene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
p-Xylene	0.2 U	0.2 U	0.2 U		0.2 U	0.2 U
m-Xylene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
o-Xylene	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Bromobenzene	_ 0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,2,3-Trichloropropane	_ 0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
I,I,I,Z-letrachloroethane	0.2 U	0.2 U	0.2 U	• • • •	0.2 U	0.2 U
n-Propy I benzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
2-Chlorotoluene	0.2 U	0.2 U	0.2 U		0.2 U	0.2 U
1-Chlorotoluene 1,3,5-Trimethylbenzene	0.2 U	0.2 U	0.2 U		0.2 U	0.2 U
,3,5-Trimethylbenzene	0.1 ป	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
ert-Butylbenzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
ert-Butylbenzene ,2,4-Trimethylbenzene	_ 0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
sec-Butylbenzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
-Isopropyltoluene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
,3-Dichlorobenzene	O.1_U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
,4-Dichlorobenzene	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
sec-Butylbenzene D-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene	O.1_U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 Ū
n-Butylbenzene 1,2-Dibromo-3-Chloropropane 1,2,4-Trichlorobenzene	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
,2-Dibromo-3-Chloropropane	_ 0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
,2,4-Trichlorobenzene	O.2 U	0.2 · U	0.2 U	0.2 U	0.2 U	0.2 U
IPEACH HIT WOILLANDEN	11 (4 1)	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
laphthalene	0.2 U	0.2 U	0.2 U	•••	0.2 U	0.2 U
,2,3-Trichlorobenzene	O.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
laphthalene ,2,3-Trichlorobenzene :is-1,3-Dichloropropene :rans-1,3-Dichloropropene	O.1_U	0.1 U	0.1 U	-	0.1 U	0.1 U
rans-1,3-Dichloropropene	0.3 U	0.3 U	0.3 U		0.3 U	0.3 U
cetone arbon Disulfide	_ 2 U	2 U	2 U		2 U	2 U
Carbon Disulfide	2	2	0.1 U		0.1 U	0.1 U
-Butanone	_ 0.3 U	0.3 U	0.3 U		0.3 U	0.3 U
/inyl Acetate	_ 0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
?-Butanone /inyl Acetate -Methyl-2-pentanone	_ 0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
2 - Hexanone	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
?-Hexanone ?-Chloroethylvinylether	0.2 U	0.2 U	0.2 U		0.2 U	0.2 U
*- Outside of EPA CLP QC limits.	_				-	- -

RFW Batch Number: 92036	G478	Client:	ULA I Ekco	ILES BY GC/MS, Houseware	METHOD 52	4 t Wor	.IST <u>rk Order; 2994-</u>	Report Date: 02-03-0000	04/15/92 11:2 Page: 2
	Cust ID:	I-13-GW	!	I-13-GW	I-14-6W	ı	Trip Blank A	Trip Blank B	
Sample	RFW#:	006 MS	,	006 MSD	007	,	008	009	010
Information	Matrix:	WATER		WATER	WATER		WATER	WATER	WATER
	D.F.:	1.0	0	1.00	1.0	n	1.00	1.00	1.00
	Units:	UG/L		UG/L	UG/L		UG/L	UG/L	UG/L
	iorobenzene		%	102 %	98	%		96 %	
)ichlorodifluoromethane)	0.3	-f1- U	fl 0.3 U	0.3	11=1 U	0.3 U	0.3 U	0.3 U
hloromethane		0.2	Ŭ	0.2 U	0.2	Ŭ	0.2 U	0.2 U	0.2
Bromomethane		0.1	Ū	0.1 U	0.1	Ŭ	0.1 U	0.1 U	0.1
/inyl Chloride		5	-	5	0.2	_	0.2 U	0.2 U	2.7
Chloroethane		0.1	U	0.1 U	0.1	Ŭ	0.1 U	0.1 U	0. Ī U
Bromomethane /inyl Chloride Chloroethane Methylene Chloride Irichlorofluoromethane 1.1-Dichloroethene		7		2	6	-	6	6	ž
richlorofluoromethane_		1		1	0.8		0.1 U	0.1 U	0.8
1,1-Dichloroethene 1,1-Dichloroethane :is-1,2-Dichloroethene		215 *	%	202 * %	0.1	U	0.1 U	0.1 U	0.1 1
,1-Dichloroethane		0.1	U	0.1 U	0.1	U	0.1 U	0.1 U	0.1
is-1,2-Dichloroethene		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	
2,2-Dichloropropane		0.1	U	0.1 U	0.1	U	0.1 U	0.1 U	0.1 l
2,2-Dichloropropane trans-1,2-Dichloroethen	ie	0.1	U	0.1 U	0.1	U	0.1 U	0.1 U	0.1 l
Chloroform		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 l
Bromochloromethane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 L
l,2-Dichloroethane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 l
l,1,1-Trichloroethane		0.2	U	0.2 · U	0.2	U	0.2 U	0.2 U	3
Carbon Tetrachloride		0.1	U	0.1 U	0.1	U	0.1 U	0.1 U	0.1 U
l,1-Dichloropropene		0.1	U	0.1 U	0.1	U	0.1 U	0.1 U	0.1 U
Chloroform Bromochloromethane 1,2-Dichloroethane 1,1,1-Trichloroethane Carbon Tetrachloride 1,1-Dichloropropene Bromodichloromethane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 (
Olbromomethane		0.2		0.2 U	0.2	U	0.2 U	0.2 U	0.2 L
Oibromomethane 1,2-Dichloropropane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2
Trichloroethene		274 *		241 * %	0.3	U	0.3 U	0.3 U	
Dibromochloromethane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2
,z-Dibromoethane		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2
1,2-Dibromoethane 1,1,2-Trichloroethane Benzene		0.3	U	0.3 U	0.3	U	0.3 U	0.3 U	0.3 (
Benzene	· · ·	162 *		160 * %	0.2	U	0.2 U	0.2 U	0.2
i,3-Ulchioropropane		U. Z	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 l
Bromoform Tetrachloroethene 1,1,2,2-Tetrachloroetha		0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2
letrachloroethene	····	0.1		0.1 U	0.1	U	0.1 U	0.1 U	0.1 (
1,1,2,2-letrachloroetha	ine	0.2	U	0.2 U	0.2	U	0.2 U	0.2 U	0.2 l
Toluene		141 *	%	139 * %	0.1	U	0.1 U	0.1 U	0.1 (
*- Outside of EPA CLP (∤C limits.								

RFW Batch Number: 9203G478	Client: Ekco	Houseware	Work	Order: 2994-0	2-03-0000	Page: 1b
Cust ID:	1-8D-GW	I -9-GW	I-10-GW	I-11-GW	I-12-GW	I-13-GW
RFW#:	001	002	003	004	005	006
Chlorobenzene	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	
1,1,1,2-Tetrachloroethane	·	0.2 U	0.2 U	0.2 U	0.2 U	0.2 Ŭ
I,I,I,Z-Tetrachloroethane Ethylbenzene Styrene p-Xylene m-Xylene o-Xylene Bromobenzene 1,2,3-Trichloropropane Isopropylbenzene n-Propylbenzene 2-Chlorotoluene 4-Chlorotoluene	O.1_U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
Styrene	_ 0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
p-Xylene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
m-Xylene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
o-Xylene	O.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Bromobenzene	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,2,3-Trichloropropane	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
Isopropylbenzene	O.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
n-Propylbenzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
2-Chlorotoluene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
4-Chlorotoluene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,3,5-Trimethylbenzene	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
tert-Butylbenzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
4-Chlorotoluene 1,3,5-Trimethylbenzene tert-Butylbenzene 1,2,4-Trimethylbenzene	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
sec-Butylbenzene p-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene 1,2-Dichlorobenzene n-Butylbenzene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
p-Isopropyltoluene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
1,3-Dichlorobenzene	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,4-Dichlorobenzene	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
1,2-Dichlorobenzene	0.1 U	0.1 U	0.1 U	0.1 Ū	0.1 U	0.1 U
1,2-Dichlorobenzene n-Butylbenzene 1,2-Dibromo-3-Chloropropane 1,2,4-Trichlorobenzene Hexachlorobutadiene Naphthalene	O.1_U	0.1 U	0.1 U	0.1 U	0.1 U	0.1 U
1,2-Dibromo-3-Chloropropane	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
1,2,4-Trichlorobenzene	0.2 บ	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
Hexachlorobutadiene	- 0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
Naphthalene	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 Ü
1,2,3-Trichlorobenzene	0.1 บ	0.1 U	0.1 U	0.1 Ü	0.1 U	0.1 U
cis-1,3-Dichloropropene	- 0.1 U	0.1 U	0.1 Ū	0.1 U	0.1 U	0.1 U
trans-1,3-Dichloropropene	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Acetone	- 2 U	2 U	2 U	2 U	2 Ū	ŽŪ
Carbon Disulfide	0.1 U	0.1 U	0.1 U	8	ī	ī
2-Butanone	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Vinyl Acetate	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U	0.3 U
Naphthalene 1,2,3-TrichTorobenzene cis-1,3-DichToropropene trans-1,3-DichToropropene Acetone Carbon Disulfide 2-Butanone Vinyl Acetate 4-Methyl-2-pentanone 2-Hexanone 2-Chloroethylvinylether *= Outside of FPA CIP OC limits	- 0.4 U	0.4 U	0.4 U	0.4 U	0.4 U	0.4 U
2-Hexanone	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U	0.5 U
2-Chloroethylvinylether	- 0.2 U	0.2 U	0.2 U	0.2 U	0.2 U	0.2 U
*- Outside of EPA CLP QC limits.	_			2.2	- • •	

RFW Batch Number: 9203647	R	Client:	VOLAT Ekco	ILÉS BY GO Houseware	/MS,	METHOD 52	24 LI Wark	ST Order: 29	 9 4 _0:	Report Dat	e: 0	4/15/92 11 Page:	:21
													•
'	Cust ID:	I - 80 - 61	1	1-9-GN	ĺ	1-10-61	•	I-11-6)	•	1-12-GM	l	1-13-64	•
Sample	RFW#:	001	l	002	!	003	•	004	ļ	005	I	006	
Information	Matrix:	WATER		WATER		WATER		WATER		WATER		WATER	
	D.F.:	1.0		1.0		1.0	00	1.0	0	1.0	0	1.0)()
	Units:	UG/L	-	UG/L		UG/L	_	UG/L	•	UG/L		UG/L	-
Bromofluore	obenzene	105	% £1	102	% -fl=:	98	*	99	% f1	102	X	103	7
Dichlorodifluoromethane				0.3	U	0.3	71-	0.3	'=T =: U	0.3	≖f1=: U	0.3	•=f1 U
chloromethane		0.2	υ	0.2	V		Ü		Ű	0.2	Ũ	0.2	Ŭ
SPOMOMPENANCE		D I	U	0.1	U	0.1	Ŭ	0.1	Ũ	0.1	Ū	0.1	Ŭ
inyl Chloride_		0.2	U	7		0.2	Ŭ	9	-	0.2	Ū	3	•
hloroethane		0.1	U	0.1	U	0.1	Ŭ	0.1	ี่ย	0.1	Ū	0.1	U
hloroethane lethylene Chloride richlorofluoromethane		5		5		9		5		4		4	
richlorofluoromethane		0.1	U	1		0.1	U	1		0.8		0.7	
,i-Dichloroethene		0.1	U	0.1	U	0.1	Ü	0.1	V	0.1	U	0.1	U
		0.1	U	0.1	U	0.1	U	1.0	U	0.1	U	0.1	Ü
,1-Dichloroethane is-1,2-Dichloroethene		0.2	U	3		0.2	U	0.2	U	0.2	U	0.2	Ū
			v	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
rans-1,2-Dichloroethene_		0.1	ช	0.1	U	0.1	U	0.1	U	0.1	U	0.1	Ü
hloroform		0.2	U	0.2	U	0.2	Ū	0.2	U	0.2	U	0.2	Ū
romochloromethane		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	V
2-Dichlornethane		0.2	U	0.2	u	0.2	U	0.2	U	0.2	U	0.2	Ü
.l.I-Trichloroethane		0.2	U	0.2	IJ	0.2	U	0.2	U	0.2	U	0.2	V
,arbon letrachioride		0.1	U	0.1	U	0.1	Ü	0.1	U	0.1	U	0.1	U
.1-Dichiaropropene		0.1	U	0.1	U	0.1	U	0.1	U	0.1	V	0.1	U
Bromodichloromethane		0.2	Ü	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
ibromomethane		0.2	U	0.2	V	0.2	U	0.2	U	0.2	U	0.2	U
,2-Dichloropropane		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
rich inroet kene			U	0.3	U	0.3	U	0.3	U	0.3	U	0.3	U
1bromochloromethane		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
,2-Dibromoethane		0.2	V	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
,2-Dibromoethane ,1,2-Trichloroethane		0.3	U	0.3	u	0.3	U	0.3	U	0.3	U	0.3	IJ
		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
,3-Dichloropropane		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
romoform		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
[etrach]oroethene		0.1	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
letrachloroethene 1,1,2,2-Tetrachloroethane		0.2	U	0.2	U	0.2	U	0.2	U	0.2	U	0.2	U
oluene		0.1	U	0.1	U	0.1	U	0.1	U	0.1	U	0.1	U
- Outside of EPA CLP QC 1	imits.												

RFW Batch Number: 9203G477		Client:	Eko	o Houseware	1		Wor	k Order: 29	94-	02-03-0000		Page: 6b
Cu	st ID:	VBLK BS		VBLK		VBLK BS		VBLK	_	VBLK BS		
	RFW#:	92GVE087-I	MB1	92GVE097-M	B 1	92GVE097-1	1B1	92GVE099-M	B 1	92GVE099-M	B 1	
Toluene		82	%	i	U	95	7	1	U	92	76	
Chlorobenzene		84	%	1	U	96	%	1	U	99	%	
Ethylbenzene		1	U	1	U	1	U	1	U	1	U	
Styrene		- 1	U	1	U	1	U	1	U	1	U	
Xylene (total)		- i	U	1	U	1	U	1	U	1	U	
*= Outside of EPA CLP QC 11	mits.	_										

RFW Batch Number: 920	36477		٧	ton, Inc (OLATILES BY (O Houseware	GC/	MS, HSL LI	ST	ratories <u>rk Order: 299</u> 4	F 1-02	Report Date: 2-03-0000	: (04/20/92 10:51 Page: 6a
	Cust ID:	VBLK BS		YBLK		VBLK BS		VBLK		VBLK BS		
Sample Information	Matrix:	92GVE087-MB WATER		92GVE097-MB1 WATER		92GVE097-M WATER	IB 1	92GVE099-MB1 WATER	L 9	92GVE099-MB1 WATER	1	
	D.F.: Units:	1.00 ug/L		1.00 ug/L		1.0 ug/l		1.00 ug/L		1.00 ug/L		
Surrogate Bromof	Toluene-d8 luorobenzene	96 90	% %		X X	101 101	% %		<u>.</u> L	94 x 102 x	X X	
Recovery 1,2-Dichl	oroethane-d4	76	%	92 %	K.	93	%	92 2	6	97 %	X.	
							•	•			-	f1
D.,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,		_ 2	U	2 U 2 U		2	U	2 L		2 4		
Vinyl Chloride		- 2	U	2 U		2	U	2 U 1 U		2 4		
Chloroethane		- 1 9	U	2 1	-	2	U	2 1	•	1 U 2 U	_	
Chloroethane Methylene Chloride		- 1	ü	1 1	_	1	Ü	1 1		2 U 1 U	_	
			Ŭ	i	•	5	В	1 (,	4 6	-	
		_	Ŭ	iu	1	1	Ü	i i		1 1		
1,1-Dichloroethene			%	iù	-	110	%	ii	-		X	
1.1-Dichioroethane		1	ũ	iŭ		110	ũ	ii		102		
1,2-Dichloroethene (to	ntall	- i	Ŭ	iŭ	-	i	Ŭ	iù		ii		
Chloroform		1	Ŭ	iŭ	_	i	Ŭ	iì	_	ii	_	
1,2-Dichloroethane		· i	Ŭ	iŭ	-	i	ŭ	ii	-	- •	U	
2-Butanone		- 2	Ŭ	2 0	-	ż	Ŭ	2 1	•		נו נו	
Z-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride		~ ī	Ŭ	ĩũ	_	້ ເ	ŭ	1 1		ii	_	
Carbon Tetrachloride		- i :	Ŭ	i ū	_	i	Ü	iù	-	iì	_	
vinvi Acetate		7	Ū	Ž Ū	j	Ž	Ŭ	żi		Ži	_	
Bromodichloromethane		- i i	Ū	īŪ	j	ī	Ū	īū		ı i i	_	
I.Z-UICHIOCODCODANE			U	1 0	j	ī	Ū	i	j	ii	_	
cis-1,3-Dichloroprope	ne	1	U	1 U	J	ī	Ũ	ii	j	īi	_	
Trichloroethene		111 9	%	1 U	J	102	%	ii	j	130 * %	X.	
Dibromochloromethane		1	U	1 8	J	1	Ü	i i	j	1 (
1,1,2-Trichloroethane		1 (U	1 U	J	ī	U	i i	j	i i	U	
Renzene		97 9	%	1 U	J	97	%	1 1	J	133 * 7	%	
Trans-1,3-Dichloropro	pene] 1 1	U	1 U	J	1	U	1 1	J	1 (U	
Bromotorm		1 1	U	1 U	_	1	U	1 1		1 (_	
4-Methyl-2-pentanone_		2	U	2 U		2	U	2 (2 l	U	
2-Hexanone		_ 2	U	2 U	_	2	U	2 l		2 l	U	
Tetrachloroethene			U	1 U		1	U	1 (1 (U	
1,1,2,2-Tetrachloroet		1	U	1 0	J	1	U	1 1	J	1 (U	

1,1,2,2-Tetrachloroethane
*- Outside of EPA CLP QC limits.

RFW Batch Number: 9203G47				o Houseware			Wor	k Order: 29	94-	02-03-0000		Page: 5b	<u>.</u>
	Cust ID:	Trip Blank	A	Trip Blank	Ā	Trip Blank	B	VBLK		VBLK BS		VBLK	
	RFW#:	012	2	012 REPR		013		92GVE086-M	B1	92GVE086-M	B1	92GVE087-MB1	
Toluene		1	U	1	U	1	Ü	1	U	105	%	1 U	_
Chlorobenzene		1	U	1	U	1	U	1	U	104	%	1 U	
Ethylbenzene		1	U	1	U	1	U	1	U	2	U	1 Ú	
Styrene		1	U	1	U	1	U	1	U	2	U	1 U	
Xylene (total)		1	U	1	U	1	U	1	U	2	U	1 U	
*= Outside of EPA CLP QC	limits.	-											

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Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS, HSL LIST
Client: Ekco Houseware Work Order: IST Report Date: 04/20/92 10:51 Work Order: 2994-02-03-0000 Page: 5a RFW Batch Number: 9203G477 Page: 5a

	Cust ID:	Trip Blank A	Trip Blank A	Trip Blank B	VBLK	VBLK BS	VBLK
Sample Information	RFW#: Matrix: D.F.: Units:	012 WATER 1.00 ug/L	012 WATER 1.00 ug/L	013 WATER 1.00 ug/L	92GVE086-MB1 WATER 1.00 ug/L	92GVEO86-MB1 WATER 2.00 ug/L	92GVE087-MB1 WATER 1.00 ug/L
	Toluene-d8	97 %	REPREP 107 %	98 %	100 %	103 %	
Surrogate Recovery	Bromofluorobenzene 1,2-Dichloroethane-d4	92 % 121 * %	93 % 94 %	90 % 80 %	101 % 106 %	97 % 70 * %	97 % 101 % 93 %
Chloromethan)6	2 U	2 U	f1 2 U	2 U	4 U	f1 2 U
Bromomethane		2 U	Ž Ū	ŽŪ	ŽŪ	4 Ŭ	2 U
Vinyl Chlori		1 U	1 Ŭ	Î Ū	ĨŪ	2 Ŭ	īŬ
Chloroethane		2 U	2 U	2 U	2 U	4 Ü	Ž Ū
Methylene Ch	loride	14	4	4	ĨŪ	2 Ŭ	ĭŭ
Acetone		6	5 B	5	Ž U	à Ū	Ž Ū
Carbon Disul		1 U	1 U	1 U	ĨŪ	2 U	ĩŬ
1,1-Dichloro		1 U	1 U	1 U	1 U	122 %	īŪ
1,1-Dichloro		1 U	1 U	1 U	1 U	2 Ü	ĬŪ
	ethene (total)	1 U	1 U	1 U	1 U	2 U	i Ū
Chloroform		1 U	1 U	1 U	1 U	2 U	i Ū
1,2-Dichloro	ethane	1 U	1 U	1 U	1 U	2 U	ÌŪ
2-Butanone		2 U	2 U	2 U	2 U	4 U	2 U
1,1,1-Trich1	oroethane	ן ו	1 U	1 U	1 U	2 U	1 U
Carbon Tetra	chloride	1 U	1 U	1 U	1 U	2 U	1 U
Vinyl Acetat	e	2 U	2 U	2 U	2 U	4 U	2 U
Bromodichlor		1 U	1 U	1 U	· 1 U	' 2 U	1 U
1,2-Dichloro		1 U	1 U	1 U	1 U	2 U	1 U
cis-1,3-Dich		1 U	1 U	1 U	1 U	2 U	1 U
Trichloroeth		. 3	4	2	1 U	114 %	1 U
Dibromochlor		. 1 U	1 U	1 U	1 U	2 U	1 U
1,1,2-Trich1	oroethane	1 U	1 U	1 U	1 U	2 U	1 U
Benzene		1 U	1 U	1 U	1 U	107 %	1 U
Trans-1,3-Di	chloropropene		1 U	1 U	1 U	2 U	1 U
Bromoform_		. 1 U	1 U	1 U	1 U	2 U	1 U
	entanone	2 U	2 U	2 U	2 U	4 U	2 บ
2-Hexanone		2 U	2 U	2 U	2 U	4 U	2 U
Tetrachloroe		1 U	1 U	1 U	1 U	2 U	1 U
	achloroethane of EPA CLP QC Timits.	1 U	1 U	1 U	1 U	2 U	1 U

RFW Batch Number: 9203G477	Client: Ekc	o Houseware	Work	Order: 2994-0	2-03-0000	Page: 4b
Cust II	D: W-10-GW	W-10-GW	S-7-GW	S-7-GW	S-7-GW	S-7-FB
RFW	#: 009 DL	009 DL	010	010 MS	010 MSD	011
Toluene	NĀ	NA NA	1 0	82 %	84 %	<u>i u</u>
Chlorobenzene	NA NA	NA	ĨŪ	84 %	85 %	ìŪ
Ethylbenzene	NA NA	NA	1 U	1 U	1 U	1 U
Styrene	NA NA	NA	1 U	1 U	1 U	1 U
Xylene (total)	NA	NA	Î Û	ĨŪ	i Ü	ĪŪ
*= Outside of FPA CLP OC limits						

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Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS. HSL LIST

RFW Batch Number: 9203647	<u> 17 </u>	<u>Client:</u>	Ekco	<u>Houseware</u>			Work	<u> Order: 29</u>	94-0	Report Dat 12-03-0000		Page:	
	Cust ID:	W-10-6W	ı	W-10-GW		S-7-GI	1	S-7-GN	İ	S-7-GW		S-7-F	В
Sample Information	RFW#: Matrix: D.F.: Units:	009 DL WATER 20. ug/l	.0	009 DL WATER 20 ug/L	0	010 WATER 1.0 ug/l	00	O10 MS WATER 1.0 ug/l	0	010 MSD WATER 1.0 ug/L	0	01 WATER 1.1 ug/	00
Surrogate Bromofluor Recovery 1,2-Dichloroe	thane-d4	101 102 99	% %	101 105 101	% %	101 91 97	% %	97 95 78	% % %	92 95 130 *		103 90 84	% %
A1 11		NA		NA	-f1	2	f1-: U	2	-†1- V	·2	-f1-: U	2	
		NA		NA		Ž	Ŭ	2	Ŭ	ž	Ŭ	ž	
		NA		NA NA		ĩ	Ŭ	1	ŭ	1	Ŭ	i	ŭ
		NA		NA NA		;	Ŭ	ż	Ŭ	ż	Ŭ	ż	Ŭ
Chloroethane Methylene Chloride		NA		NA NA		Ä	ŭ	3	•	Ä	•	Ä	U
Acetone		NA NA		NA NA		ż	ΰ	2	U	,	U	,	U
Carbon Disulfide		NA NA		NA		ĩ	Ŭ	1	Ŭ	1	Ŭ	í	ย
1,1-Dichloroethene		NA NA		NA		i	Ŭ	95	%	92	%	i	Ŭ
1,1-Dichloroethane		100		NA NA		P73.		2	~	2	~	i	ŭ
1,2-Dichloroethene (total		200		NA NA		7	U	1	U	1	U	;	Ŭ
Chloroform	/	NA NA		NA NA		i	Ŭ	i	Ŭ	i	ŭ	i	ŭ
1,2-Dichloroethane		NA NA		NA NA		1	Ü	1	Ŭ	;	ŭ	1	ŭ
2-Butanone		NA NA		NA NA		1	ŭ	1	Ŭ	2	Ŭ	2	Ü
1,1,1-Trichloroethane		WA	E	1600		F7 '		£	U	7	U	,	Ü
Carbon Tetrachloride		NA	Ľ	NA .		`,	U	1	U	1	u	•	Ü
Vinyl Acetate		NA NA		NA NA		1	Ü	1	Ü	1	ŭ	1	Ü
Promodichloromethane		NA NA		NA NA		. 1	Ü	: 1	Ü	; 1	Ü	1	U
		NA NA		NA NA		1	Ü	1	ŭ	1	Ü	1	U
1,2-Dichloropropane cis-1,3-Dichloropropene		NA NA		NA NA		1	Ü	1	Ü	i	Ü	1	u
Trichloroethene		IVA	Ε	790		*9	٠	85	₹.	74	%	1	J
Dibromochloromethane		NA	C	NA		" y	U	69 1	ũ	71	Ű	1	U
1,1,2-Trichloroethane		NA NA		NA NA		1	Ü	, , , , , , , , , , , , , , , , , , ,	Ü	1	Ü	1	Ü
1,1,2-111CHTOFOECHWHE Benzene		NA NA		NA NA		1	Ü	101	%	149	_	1	Ü
Trans-1,3-Dichloropropene		NA NA		NA NA		1	IJ	101	ũ	149	ΰ	1	Ü
		NA NA		NA NA		1	Ü	1	Ü	i	Ŭ	1	U
Bromoform 4-Methyl-2-pentanone		NA NA		NA NA		2	_	1	Ü	2	Ŭ	1	U
T-methyl-c-pentanone		NA NA		NA NA		2	IJ	2	Ŭ	2	Ŭ	2	; U
2-Hexanone				NA NA		1	Ü	1	Ŭ	1	Ü	1	U
		NA NA		NA NA		1	U	1	Ü	1	Ü	1	U I II
1,1,2,2-Tetrachloroethane *= Outside of EPA CLP QC	•	NA		MA		1	U	1	U	1	V	1	U

FW Batch Number: 9203G477		Client: Ekc				2-03-0000	Page: 3b				
	Cust ID:	I-2-GW-Dup	I-5-GW		I-6-GW		W-1-GW		W-1-GW	W-10-6W	ı
	RFW#:	005 DL	006		007		800		008 DL	009)
To l uene	- · · · · · · · · · · · · · · · · · · ·	NA	i	U	i	Ū	i	Ū	NA NA	0.6	_j
Chlorobenzene		- NA	1	U	1	U	1	U	NA	1	U
thylbenzene		- NA	1	U	1	U	1	U	NA	1	U
Styrene		- NA	1	U	1	U	1	U	NA	1	U
(ylene (total)		- NA	1	U	1	U	1	U	NA	i	U

Roy F. Weston, Inc. - Gulf Coast Laboratories

RFW Batch Number: 9203G477	Client:	V Ekc	OLATILES BY <u>o Houseware</u>	GC,	/MS, HS			k Order: 29	94-	Report Date: 0 -02-03-0000	04/20/92 10 Page:	0:51 3a
Cust ID:	I-2-GW-Dup)	I-5-GW		I-	6-GW	1	W-1-6k	1	W-1-GW	M-10-6K	
Sample RFW#: Information Matrix: D.F.: Units:	WATER 50.	0	006 WATER 1.00 ug/L	0		007 TER 1.0 ug/L	0	008 WATER 1.0 ug/l	00	008 DL WATER 5.00 ug/L	009 WATER 1.0 ug/L	00
Toluene-d8	98	%	101	*		05	*	97	%	99 %	98	
Surrogate Bromofluorobenzene Recovery 1,2-Dichloroethane-d4	102 97	% %	100 78	% %	1	98 00	% %	90 80	% %	100 % 92 %	90 79	% % %
Chloromethane	NA NA	- † [·	2	-f1- U		2	-f1- U	2	f1-י ט	f1- NA		fi U
Bromomethane	- NA		Ž	Ū		Ž	Ū	2	Ŭ	NA NA	2	Ü
Vinyl Chloride	NA		ī	U		Ī	Ū	ī	Ū	NA	3	v
LNIOTOEINANE	- NA		2	U		2	Ŭ	2	Ŭ	NA	2	U
Methylene Chloride	- NA		3	U		4	_	3	Ū	NA	3	•
Acetone	_ NA		2	Ŭ		2	U	2	U	NA NA	ž	U
Carbon Disulfide	NA		1	U		Ĩ	Ũ	ī	Ŭ	NA	ĩ	Ŭ
1,1-Dichloroethene	_ NA		1	U		Ī	Ū	8	•	NA	2 i	•
1,1-Dichloroethane	920	•	18	•		1	Ū	_	Ε	70	~•	F
1,2-Dichloroethene (total)			1	U		1	Ū	12	_	NA		Ē
Chloroform	- NA		1	U		ĺ	U	Ĭ	U	NA	1	ũ
1,2-Dichloroethane	_ NA		1	U		1	U	Ĭ	Ū	NA	i	Ū
2-Butanone	- NA		2	U		2	Ŭ	Ž	Ū	NA	į	Ŭ
1,1,1-Trichloroethane	- NA		1.	U		1	U		Ē	37	•	F
Lardon letrachioride	- NA		1	U		1	Ū	1	Ũ	NA	1	ũ
Vinyl Acetate	NA		2	U		2	U	2	U	NA	ž	Ū
Bromodichloromethane	- NA		1	U		1	U	Ī	U	NÀ	i	Ū
1,2-Dichloropropane	NA		1	U		1	U	Ī	U	NA	i	Ū
cis-1,3-Dichloropropene	NA NA		1	U		1	U	1	U	NA	ĺ	U
Trichloroethen e	520	•	1	•		1			Ε	140		E
Dibromochloromethane	NA NA		1	U		1	U	1	U	NA	1	Ü
1,1,2-Trichloroethane	_ NA		1	U		1	U	1	U	NA	ĺ	U
Benzene	NA		1	U		1	U	1	U	NA	1	U
Trans-1,3-Dichloropropene			1	U		1	U	1	U	NA	1	U
Bromoform	NA		1	U		1	U	1	U	NA	1	U
4-Methy1-2-pentanone	_ NA		2	U		2	U	2	U	NA	2	Ū
2-Hexanone	_ NA		2	U		2	U	2	U	NA	2	Ü
Tetrachloroethene	_ NA		1	U		1	U	1	U	NA	ĩ	Ū
1,1,2,2-Tetrachloroethane	_ NA		1	U		1	U	1	U	NA	Ĭ	Ŭ
*= Outside of EPA CLP QC limits.	_										-	-

RFW Batch Number: 9	Cust ID:	Client: Ekco D-4-30-GW	1-2-GW		I-2-GW	<u> Order: 2994</u> I-2-GW	1-2-GW-Dup	i	Page: -2-GW-Dup
	RFW#:	003 DL	004		004 DL	004 DL	005		005 DL
Toluene		NÁ	i	Ü	NA	NA	<u> </u>	Ü	NA
Chlorobenzene		NA	1	U	NA	NA	ĺ	U	NA
Ethylbenzene		. NA	1	U	NA	NA	Ī	U	NA
Styrene		NA	1	U	NA	NA	ī	U	NA
Xylene (total)		NA	ī	U	NA	NA	i	Ū	NA
*= Outside of EPA C	LP OC limits.	•	_				_	•	

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Roy F. Weston, Inc. - Gulf Coast Laboratories

RFW Batch Number: 92	036477	Client:	V(Ekc	DLATILES BY Houseware	GC/	MS, HSL L	IST Work	Order: 2994-	Report Date: 02-03-0000	04/20/92 10:5 Page: 2
	Cust ID:	D-4-30-6	W	1-2-GW		I-2-61		1-2-GW	I-2-GW-Dup	I-2-GW-Dup
Sample	RFW#:	003 D	L	004		004 DI	L	004 DL	005	005 DL
Information	Matrix:	WATER		WATER		WATER	_	WATER	WATER	WATER
	D.F.:	40	00	1.0	0	5.0		50.0	1.00	5.00
	Units:	ug/	L	ug/L		ug/l	L	ug/L	ug/L	ug/L
Summara to Durant	Toluene-d8	95	%	94	*	103	%	102 %	96 %	99 %
Surrogate Bromot Recovery 1,2-Dichi	fluorobenzene loroethane-d4	103 99	% %	91 78	% %	97 94	% %	101 % 97 %	93 % 75 * %	97 % 108 %
	****				-f1-		f1-	f1	f1	f
	····			2	ñ	NA		NA	2 U	NA
Bromomethane Vinyl Chloride		NA NA		2	ñ	NA		NA	2 U	NA.
Chloroethane	 -	. NA NA		ခဋ	E	,.59		NA	E	#P3
Chloroethane Methylene Chloride		. NA NA		•	u	NA NA		NA	PIG	NA
				3	U	NA NA		NA	4 U	NA
Carbon Disulfide		. NA NA		2	Ü	na NA		na Na	2 U	NA
				1	E	770	٠,		า กั	NA ** 6 6
1 1 Dichlementhame	······································	AI A			ב	,70	Ε	NA 980 ° '	į.	. 60
1,2-Dichloroethene (i	ntall	NA NA			Ē		Ē	760	ž r	E E
Chloroform		. NA		3	ũ	NA	C	NA NA	1 0	NA E
1,2-Dichloroethane		NA NA		i	Ŭ	NA NA		NA NA	921 *** ·	NA NA
A D. A		• • • •		Ž	Ŭ	NA		NA		NA NA
1.1.1-Trichloroethane		NA		-	Ĕ	7 THE 7.	• •	NA	2 U 725 '	NA NA
Carbon Tetrachloride		NA		1	Ū	NA NA		NA	1 U	NA NA
Vinvì Acetate		NΔ		Ž	Ŭ	NA		NA	ŽŬ	NA NA
Bromodichloromethane		NA		ī	Ū	NA		NA	iiū	NA
1,2-Dichloropropane ~		NA		Ĭ	Ū	NA		NA	īŪ	NA
cis-1,3-Dichloroprope	ene	NA		ī	U	NA		NA	iŭ	NA
Trichioroethene	***************************************	725000			E		E	490	Ē	E
Dibromochloromethane		NA		1	U	NA		NA	1 Ũ	NA ~
1,1,2-Trichloroethane		NA		1	U	NA		NA	i U	NA
Benzene		NA		1	U	NA		NA	1 U	NA
Trans-1,3-Dichloropro	pene	NA		1	U	NA		NA	1 U	NA
N				1	U	NA		NA	1 U	, NA
Bromoform 4-Methyl-2-pentanone_		NA		2	U	NA		NA	S A	NA
2-Hexanone		NA		2	U	NA		NA	2 U	NA
Tetrachloroethene		NA		1	U	NA		NA	1 U	NA
1,1,2,2-Tetrachloroet		NA		1	U	NA		NA	1 U	NA
*- Outside of EPA CLF	QC limits.									

Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS, HSL LIST

RFW Batch Number: 910	96745		ston, Inc G VOLATILES BY G co Houseware	C/MS, HSL LIST	ratories rk_Order: 2994-	Report Date:	10/15/91 14:05 Page: 3a	
na baten namet, 520	Cust ID:	D-4 -30	VBLK	VBLK BS	VBLK	VBLK BS	VBLK	
Sample	RFW#:	005 DL	91GVB342-MB1	91GVB342-MB1	91GVB343-MB1	91GVB343-MB1	91GVB344-MB1	
Information	Matrix:	WATER	WATER	WATER	WATER	WATER	WATER	
	D.F.:	500	1.00	1.00	1.00	1.00	1.00	
	Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L	
	Toluene-d8	104 %				105 %	99 %	
	luorobenzene	104 %		107 %		118 * %	100 %	
Recovery 1,2-Dichlo	oroethane-d4	104 %	93 %	104 %		112 %	95 %	
		NA NA	10 U	10 U	10 U	10 U	10 U	
		NA	10 U	10 U	10 U	10 U	10 0	
		NA	10 U	10 U	10 U	10 U	10 U	
-11		NA	10 U	10 U	10 U	10 U	10 U	
Chloroethane Methylene Chloride		NA	5 U	5 U	3 J	15 B		
Acatana		NA	17	10 U	8 J	10 U	10 U	
		NA	5 ป	5 U	5 U	5 U	5 U	
1,1-Dichloroethene		NA	5 U	85 %	5 U	90 %	5 U	
1,1-Dichloroethane		NA	5 U	5 U	5 U	5 U	5 U	
1,2-Dichloroethene (to	otal)	NA	5 U	5 U	5 U	5 U	5 U	
Chloroform		NA	5 U	5 U	5 U	5 U	5 U	
1,2-Dichloroethane		NA	5 U	5 U	5 V	5 U	5 U	
2-Butanone		NA	10 U	10 U	10 U	10 U	10 U	
1,1,1-Trichloroethane		NA	5. U	11	5 U	5 U	5 U	
Carbon letrachloride_		NA	5 U	5 U	5 U	5 U	5 U	
Vinyl Acetate		NA	10 U	10 U	10 U	10 U	10 U	
Bromodichloromethane_	···	NA	5 U	5 U	5 U	5 U	5 U	
1,2-Dichloropropane	 -,	NA	5 U	5 U	5 U	5 U	5 U	
cis-1,3-Dichloroproper	ne	NA	5 ป	5 U	5 U	5 V	5 U	
Trichloroethene		75000	5 U	246 * %	5 U	99 %	5 U	
Dibromochloromethane_		NA	5 U	5 U	5 U	5 U	5 U	
1,1,2-Trichloroethane		NA	5 U	5 U	5 U	5 U	5 U	
Benzene		NA	5 U	96 %	5 Ų	103 %	5 V	
Trans-1,3-Dichloroprop	pene	NA	5 U	5 U	5 U	5 U	5 U	
Bromoform		NA	5. U	5 U		5 U	5 U	
4-Methyl-2-pentanone_		NA	10 U	10 U		10 U	10 U	
2-Hexanone		; NA	10 U	10 U		10 U	10 U	
Tetrachloroethene		NA	5 U	5 U		5 U	5 U	
1,1,2,2-Tetrachioroet	hane	NA	5 V	5 U	5 U	5 U	5 U	
*= Outside of EPA CLP	ŲČ limits.							

RFW Batch Number: 9109G745	<u> Client: Ek</u>	co Houseware	<u>rk Order: 2994-</u>	02-03-0000	Page: 3b	
Cust ID:	0-4-30	VBLK	VBLK BS	VBLK	VBLK BS	VBLK
RFW#:	005 DL	91GVB342-MB1	91GVB342-MB1	91GVB343-MB1	91GVB343-MB1	91GVB344-MB1
Toluene	NA	5 U	92 %	5 11	98 %	5 U
Chlorobenzene	- NA	5 Ū	94 %	5 Ŭ	94 %	5 Ŭ
Ethy1benzene	- NA	5 Ú	5 Ü	5 Ŭ	5 11	5 Ŭ
Styrene	- NA	5 Ū	5 Ü	5 Ŭ	5 Ü	5 Ü
Xylene (total)	- NA	5 Ū	5 Ü	5 Ü	5 Ŭ	5 Ŭ
2-Chloroethylvinylether *= Outside of EPA CLP QC limits.	NA	10 U	10 U	10 Ü	10 U	10 Ü

Roy F. Weston, Inc. - Gulf Coast Laboratories VOLATILES BY GC/MS, HSL LIST

RFW_Batch_Number: 9109	0G745		VO		GC/I	MS, HSL LI	ST	Order: 29	Rep 94-02-0	ort Da 3-0000	te: 10	/15/91 1 Page:	
	Cust ID:	R-5-G	!	R-5-GW		R-5-GW		D-4-30		D-4-30		D-4-3	
Sample Information	RFW#: Matrix: D.F.: Units:	004 WATER 1.0 ug/l	0	004 MS WATER 1.0 ug/L	0.	004 MSD WATER 1.0 ug/L	10	005 WATER 1.0 ug/L	0	005 DI WATER 10 ug/	.0	OOS D Water 1 ug/	00
	Toluene-d8 uorobenzene proethane-d4	100 106 103	% % %	103 111 107	% % %	103 109 106	% % %	108 108 108	% % %	103 115 113	% % %	98 100 100	% % %
Chloromethane Bromomethane Vinyl Chloride Chloroethane Methylene Chloride Acetone Carbon Disulfide 1,1-Dichloroethane 1,2-Dichloroethane 1,2-Dichloroethane 2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride Vinyl Acetate Bromodichloromethane 1,2-Dichloropropane cis-1,3-Dichloropropen	re	10 34 10 5 5 5 5 5 10 5 5 5 5 5 5 5 5 5 5 5 5		10 10 29 10 5 10 5 7 54 5 10 35 10 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5 5	טעטעע טעט %עטטע	10 10 29 10 5 10 5 89 6 53 5 10 17 5 5 5 10 5 5 5 5 5 5 5 5 5 10 5 5 5 5 5	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	10 10 10	U U U U	NA NA NA NA NA 930 1600 NA NA NA NA NA	•	NA N	

RFW Batch Number: 91096745			<u>Houseware</u>			Work	Order: 299	94-02	-03-0000	Page: 2b
Cust ID:	R-5-GI	H	R-5-GK	1	R-5-GW		D-4-30		D-4-30	D-4-30
RFW#:	004	\$	004 MS	;	004 MSD)	005		005 DL	005 DL
toluene	5	U	90	%	95	%			NA	NA
Chlorobenzene	5	U	91	%	98	%	5	U	NA	NA
Ethylbenzene	5	U	5	U	5	Ü	5	Ū	NA	NA
Styrene	5	U	5	U	5	U	5	Ü	NA	NA
Xylene (total)	5	U	5	U	5	U	5	Ū	NA	NA
2-Chloroethylvinylether	10	U	10	U	10	U	10	Ū	NA	NA
*= Outside of EPA CLP QC limits.	110	>						970		····

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Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS. HSL LIST

RFW Batch Number: 91096	\7 A P	0114.	rkaa	Houseven	. 00/11	S, HSL LI				Report Date	:: 10/		
	1/45	Cilent:	LKCO	<u>Housewar</u>	9		work C	raer: 29	<u> 14-U</u>	2-03-0000		Page:	<u> 1 a</u>
	Cust ID:	I - 7 - GV	1	I-7-GI	1	I-2-GW	1	I-2-GW		I-7-GW DUP	1 -	7-GW DUI	Þ
Sample Information	RFW#: Matrix:	001 WATER	l	001 DI WATER	L	002 WATER	!	002 DL WATER		003 WATER		003 DI	
Intormation	D.F.:	1.0	10	WATER 5.0	30	WATER 1.0	10	WATER 5.0	n	WATER 1.00	`	WATER 5.0	
	Units:	ug/l		ug/l	-	ug/L		ug/L	U	ug/L	-	ug/	
	Toluene-d8	92	7,	104	%	100	%	106	%	102	%	104	%
	ıorobenzene	112	%	109	%	108	%	110	%	104	%	108	%
Recovery 1,2-Dichlor	roethane-d4	108	%	111	%	107	%	110	%	102	%	112	%
*************					==f]==				=f]=				==[]
Chloromethane		10	Ü	NA			U	NA			U	NA	
Bromomethane Vinyl Chloride		10	U	NA		10	ָ, U	NA		10	U	NA	
Vinyl Chloride		10	V	NA		76		NA		10	U	NA	
Chloroethane Methylene Chloride		10	U	NA		10	U	NA		10	U	NA	
Methylene Chloride		. 5	U	NA		5	U	NA		5	U	NΛ	
Acetone		10	U	NA		8	$JB\mathcal{K}$	NA		10	U	NA	
Acetone Carbon Disulfide		5	U	NA		5	U	NA		5	U	NA	
1,1-Dichloroethene		5	U	NA		*85 "	1: :	NA		6 7 11		NA	
1,1-Dichloroethene 1,1-Dichloroethane 1,2-Dichloroethene (tot			Eilis		*		E 950		•		E 480	480	
1,2-Dichloroethene (tot	:al)	5	U	NA		₹130	•	NA		72	•	NA	
Chloroform		5	U	NA		5	U	NA		5	U	NA	
Chloroform 1,2-Dichloroethane		5	U	NA		5	U	NA		5	U	NA	
2-Butanone		10	U	NA		10	U	NA		10	U	NA	
1,1,1-Trichloroethane_		- 5	U	1473		725		NA		5	U	NA	
2-Butanone 1,1,1-Trichloroethane Carbon Tetrachloride		5	U ´	NA		5	U	NA		5	U	NA	
Vinyl AcetateBromodichloromethane		10	U	NA		10	U	NA		10	U	NA	
Bromodichloromethane		5	U	NA		5	U	NA		5	U	NA	
1,2-Dichloropropane		5	U	NA		5	U	NA		5	U	NA	
1,2-Dichloropropanecis-1,3-Dichloropropene	<u> </u>	5	U	NA		5	U	NA		5	U	NA	
Trichloroethene		5	U	NA			E 1/61	7 7460		*'6 · ·		NA	
Dibromochloromethane		5	U	NA		5	U	NA		5	บ	NA	
1,1,2-Trichloroethane_		5	U	NA		5	U	NA		5	U	NA	
^		5	U	NA		5	U	NA		5	U	NA	
Trans-1,3-Dichloroprope	ne	5	Ū	NA		5	Ū	NA		5	Ũ	NA	
		<u> </u>	Ŭ	NA		5	Ū	NA		5	Ū	NΑ	
Bromoform 4-Methyl-2-pentanone	·	10	Ū	NA		10	Ū	NA		10	Ū	NΛ	
2-Hexanone		: 10	Ū	NA		10	Ŭ	NA		10	Ŭ	NΑ	
Tetrachloroethene		5	Ŭ	NA		5	Ŭ	NA		5	Ŭ	HΛ	
1,1,2,2-Tetrachloroetha		5	Ŭ	NA		5	Ŭ	NA		5	Ŭ	NΛ	
*= Outside of EPA CLP (C limits.	3	-	• • • • • • • • • • • • • • • • • • • •		•	•	. 47 5		J	-		

RFW Batch Number: 9109G745	<u>Client:</u>	Ekco	Houseware		Work	Order: 2994	-02-03-0000		Page: 1b
Cust ID:	1-7-GW		I-7-GW	1-2-GW		1-2-GW	1-7-GW DUP		1-7-GW DUP
RFW#:	001		001 DL	002		002 DL	003		003 DL
Toluene	5	U	NĀ	5	U	NĀ	5	U	NA NA
Chlorobenzene	5	U	NA	5	U	NA	5	Ũ	NA
Ethylbenzene	5	U	NA	5	U	NA	5	Ŭ	NA
Styrene	5	U	NA	5	U	NA	5	Ū	NA
Xylene (total)	5	U	NA	5	U	NA	5	Ŭ	NA
2-Chloroethylvinylether	10	U	NA	10	Ū	NA	_	Ŭ	NA
*= Outside of EPA CLP QC limits.	F_1	/)		6 مارا	2		50		7

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Client: Ekco Houseware

RFW Batch Number: 9109G723

*= Outside of EPA CLP QC limits.

VOLATILES BY GC/MS, METHOD 524 LIST

Report Date: 10/11/91 11:25 CD

Work Order: 2994-02-03-0000

Page: 2a Cust ID: **VBLK VBLK BS** Sample RFW#: 91GVE242-MB1 91GVE242-MB1 Information Matrix: WATER WATER D.F.: 1.00 1.00 Units: UG/L UG/L Bromof Luorobenzene 113 114 Dichlorodifluoromethane 0.3 U 0.2 U 0.2 U Chloromethane 0.1 U Bromomethane 0.1 U Vinyl Chloride 0.2 U U Chloroethane 0.1 U Methylene Chloride Trichlorofluoromethane U U 1.1-Dichloroethene U % 1.1-Dichloroethane cis-1,2-Dichloroethene U 0.2 U 2,2-Dichloropropane · 11 trans-1,2-Dichloroethene Chloroform U **Bromochloromethane** 0.2 u 1,2-Dichloroethane 0.2 u 1.1.1-Trichloroethane U U Carbon Tetrachloride u 1.1-Dichloropropene 0.1 U Bromodichloromethane Dibromomethane U 1.2-Dichloropropane 0.2 U Trichloroethene Dibromochloromethane 0.2 U 1.2-Dibromoethane U 1,1,2-Trichloroethane U Benzene 1,3-Dichloropropane 0.2 U Bromoform 0.2 U Tetrachloroethene 0.1 U 1,1,2,2-Tetrachloroethane U 0.2 0.2 U Toluene

Page: 2b

RFW Batch Number: 9109G723 Client: Ekco Houseware Work Order: 2994-02-03-0000 Cust ID: VBLK VBLK BS RFW#: 91GVE242-MB1 91GVE242-MB1 Chlorobenzene 0.1 U 86 1,1,1,2-Tetrachlorethane 0.2 U 0.2 U Ethylbenzene____ 0.1 U 0.1 11 Styrene 0.2 U 0.2 U p-Xylene_____ 0.2 U 0.2 U m-Xylene 0.2 U 0.2 U o-Xvlene 0.3 U 0.3 U Bromobenzene U 0.1 0.1 1 1.2.3-Trichloropropane U 0.4 0.4 U Isopropylbenzene 0.2 U 0.2 U n-Propylbenzene 0.2 U 0.2 U 2-Chlorotoluene 0.2 U 0.2 U 4-Chlorotoluene 0.2 U 0.2 .U 1.3.5-Trimethylbenzene 0.1 U 0.1 U tert-Butylbenzene 0.2 U 0.2 U 1.2.4-Trimethylbenzene 0.3 U 0.3 U sec-Butvlbenzene 0.2 U 0.2 U p-Isopropyltoluene 0.2 U 0.2 U 1,3-Dichlorobenzene 0.1 U 0.1 U 1.4-Dichlorobenzene 0.3 U 0.3 U 1,2-Dichlorobenzene 0.1 U 0.1 U 0.1 U n-Butylbenzene 0.1 U 1,2-Dibromo-3-Chloropropane 0.5 U 0.5 U 1.2.4-Trichlorobenzene 0.2 U 0.2 U Hexachlorobutadiene 0.4 U 0.4 U Naphthalene 0.2 U 0.2 U 1,2,3-Trichlorobenzene U 0.1 0.1 U cis-1,3-Dichloropropene 0.1 U 0.1 U trans-1,3-Dichloropropene 0.3 U 0.3 U Acetone 2 U 2 U Carbon Disulfide 0.1 U 0.1 U 2-Butanone 0.3 U 0.3 U Vinyl Acetate 0.3 U 0.3 U 4-Methyl-2-pentanone 0.4 U 0.4 U 2-Hexanone 0.5 U 0.5 U 2-Chloroethylvinylether 0.2 U 0.2 U *= Outside of EPA CLP QC limits.

Roy F. Weston, Inc. - Gulf Coast Laboratories

VOLATILES BY GC/MS, METHOD 524 LIST Report Date: 10/11/91 11:25 RFW Batch Number: 9109G723 Client: Ekco Houseware Work Order: 2994-02-03-0000 S Page: la Cust ID: 1-13-GW R-4-GW I-4-GW I-4-GW **VBLK** VBLK BS Sample RFW#: 001 002 003 003 DL 91GVE239-MB1 91GVE239-MB1 Information Matrix: WATER WATER WATER WATER WATER WATER D.F.: 1.00 1.00 1.00 5.00 1.00 1.00 Units: UG/L UG/L UG/L UG/L UG/L UG/L Bromof luorobenzene 116 116 117 % 116 % 121 Dichlorodifluoromethane 0.3 0.3 U 0.3 U NA 0.3 U 0.3 Chloromethane 0.2 U 0.2 U 0.2 U NA 0.2 U 0.2 U Bromomethane 11 0.1 U NA 0.1 U 0.1 Vinyl Chloride 2 0.2 U NA 0.2 U Chloroethane H 0.1 · U 0.1 U NA 0.1 U Methylene Chloride В 2 BU 2 NA 2 2 Trichlorofluoromethane U U 0.1 0.1 U NA U 0.1 1,1-Dichloroethene 0.1 U U NA 1) 94 1,1-Dichloroethane 2 164 0.1 U U cis-1,2-Dichloroethene 0.2 U 0.2 u NA 0.2 U 0.2 U U 2,2-Dichloropropane Ħ NA U trans-1.2-Dichloroethene 0.1 U 0.1 U NA 0.1 U Chloroform 0.2 U 0.2 U 0.2 NA 0.2 11 Bromochloromethane H NA 0.2 U U 1,2-Dichloroethane 0.2 0.2 0.2 NA 0.2 U 1,1,1-Trichloroethane 0.2 11 Ħ NA U Carbon Tetrachloride U NA 0.1 U U 1.1-Dichloropropene u NA 0.1 U Bromodichloromethane 0.2 0.2 U U NA 0.2 U U 0.2 Dibromomethane 0.2 0.2 0.2 U NA 0.2 U U 1,2-Dichloropropane 0.2 0.2 u 0.2 NA 0.2 U Trichloroethene 0.3 NA 0.3 U 93 Dibromochloromethane 0.2 IJ 0.2 U 0.2 U NA 0.2 U 0.2 U 1,2-Dibromoethane 0.2 0.2 U NA 0.2 U 0.2 1.1.2-Trichloroethane 0.3 NA 0.3 U 0.3 U Benzene 0.2 U 0.2 U 0.2 U NA 0.2 U 100 1,3-Dichloropropane 0.2 U 0.2 IJ NA 0.2 U 0.2 U Bromoform 0.2 0.2 0.2 U NA 0.2 U 0.2 U Tetrachloroethene 0.1 U NA 0.1 U U 1,1,2,2-Tetrachloroethane 0.2 U 0.2 U 0.2 U NA 0.2 U 0.2 U Toluene 0.9 0.1 U 0.1 U NA 0.1 U 96 *= Outside of EPA CLP OC limits.

W Batch Number: 9109G723 Cust ID:	Client: Ekco I-13-GW		<u>Work</u>	<u> Order: 2994</u>		Page: 1b
cust to:	1-13-GW	R-4-6W	1-4-GW	1-4-GW	VBLK	VBLK BS
RFW#:	001	002	003	003 DL	91GVE239-MB1	91GVE239-MB1
hlorobenzene ,1,1,2-Tetrachlorethanethylbenzene tyrene -Xvlene	0.1 U	0.1 U	0.1 U	NA	0.1 U	101 %
,1,1,2-Tetrachlorethane	0.2 U	0.2 U	0.2 U	NA NA	0.2 U	0.2 U
thylbenzene	0.1 U	0.1 U	0.1 U	NA	0.1 U	0.1 U
tyrene	0.2 U	0.2 U	0.2 U	NA NA	0.1 U	0.1 U
- Xylene	0.2 U	0.2 U	0.2 U	NA	0.2 U	0.2 U
- Xylene - Xylene - Xylene	0.2 U	0.2 U	0.2 U	NA NA	0.2 U	0.2 U
- Xy lene	0.3 U	0.3 U	0.3 U	NA NA	0.3 U	0.2 U
romobenzene	0.1 U	0.1 U	0.1 U	NA	0.1 U	0.3 U
-Xylene romobenzene ,2,3-Irichloropropane sopropylbenzene -Propylbenzene -Chlorotoluene -Chlorotoluene -Chlorotoluene ,3,5-Trimethylbenzene ert-Butylbenzene ert-Butylbenzene -2,4-Trimethylbenzene ec-Butylbenzene -Isopropyltoluene ,3-Dichlorobenzene ,4-Dichlorobenzene ,2-Dichlorobenzene -Butylbenzene ,2-Dibromo-3-Chloropropane ,2,4-Trichlorobenzene exachlorobutadiene aphthalene ,2,3-Trichlorobenzene is-1,3-Dichloropropene rans-1,3-Dichloropropene	0.4 U	0.4 U	0.4 U	NA NA	0.1 U	0.1 U
sopropylbenzene	0.2 U	0.2 U	0.2 U	NA NA	0.7 U	0.2 U
- Propylbenzene	0.2 U	0.2 U	0.2 U	NA	0.2 U	0.2 U
-Chlorotoluene	0.2 U	0.2 U	0.2 Ŭ	NA	0.2 U	0.2 U
-Chlorotoluene	0.2 U	0.2 U	0.2 U	NA	0.2 U	0.2 U
,3,5-Trimethylbenzene	0.1 U	0.1 U	0.1 U	NA	0.1 U	0.1 U
ert-Butylbenzene	0.2 U	0.2 U	0.2 U	NA	0.2 U	0.2 U
.2.4-Trimethylbenzene	0.3 U	0.3 U	0.3 U	NA	0.3 U	0.3 U
ec-Butylbenzene	0.2 U	0.2 U	0.2 U	NA	0.2 U	0.3 U
-Isopropyltoluene	0.2 U	0.2 U	0.2 Ŭ	. NA	0.2 U	0.2 U
.3-Dichlorobenzene	0.1 U	0.1 U	0.1 Ŭ	NA	0.1 U	0.1 U
.4-Dichlorobenzene	0.3 U	0.3 U	0.3 U	NA NA	0.3 U	0.1 U
.2-Dichlorobenzene	0.1 U	0.1 U	0.1 U	NA NA	0.1 U	0.3 U
-Butylbenzene	0.1 U	0.1 U	0.1 U	NA NA	0.1 U	0.1 U
.2-Dibromo-3-Chloropropane	0.5 U	0.5 U	0.5 U	NA NA	0.5 U	0.1 U
.2.4-Trichlorobenzene	0.2 U	0.2 U	0.2 U	NA NA	0.3 U	0.5 U
exachlorobutadiene	0.4 U	0.4 U	0.4 U	NA NA	0.4 U	0.2 U
aphthalene	0.2 U	0.2 U	0.2 Ŭ	NA NA	0.7 U	0.7 U
.2.3-Trichlorobenzene	0.1 U	0.1 U	0.1 U	NA NA	0.1 U	0.2 U
is-1.3-Dichloropropene	0.1 U	0.1 U	0.1 U	NA NA	0.1 U	0.1 U
rans-1.3-Dichloropropene	0.3 U	0.3 U	0.3 U	NA NA	0.1 U	0.1 U
cetone	. 2 U	2 U	2 U	NA NA	2 0	2 U
cetone arbon Disulfide	0.1 0	0. Ī Ū	0.Ì Ŭ	NA NA	0.1 U	0.1 U
-Butanone	0.3 U	0.3 U	0.3 U	NA NA	0.1 U	0.1 U
invl Acetate	0.3 U	0.3 U	0.3 U	NA NA	0.3 U	0.3 U
-Methyl-2-pentanone	0.4 Ŭ	0.4 U	0.4 U	NA NA	· 0.4 U	0.3 U 0.4 U
arbon Disulfide -Butanone inyl Acetate -Methyl-2-pentanone -Hexanone	0.5 U	0.5 U	0.5 U	NA NA	0.4 U	
-Hexanone -Chloroethylvinylether	0.2 U	0.3 U	0.3 U	NA NA	0.5 U 0.2 U	0.5 U 0.2 U

RFW Batch Number: 920	126410	-	V	ton, Inc OLATILES BY O Houseware	GC/MS, HSL	Laboratories LIST	Report Date 2994-02-02-0000	: 02/25/92 14:49 Page: 5a
KIN DELCH NUMBER . 320	69717	CITALL	EKÇ	O HORZEMELE		MALK OLDER	2734-02-02-0000	
	Cust ID:	VBLK BS						
Sample		92GYE024-1	181					
Information	Matrix:	WATER						
	D.F.:	1.9						
	Units:	ug/l	-					
	Toluene-d8	100	*					
	luorobenzene	105	%					
	oroethane-d4	102	%		£1	61		f]f]
Chloromethane	***********	2	U 7 P		.,		224 (
Bronomethane	······································	- 2	Ŭ					
Vinyl Chloride		- î	ŭ					
Chloroethane		- į	ŭ					
Methylene Chloride			Ū					
Acetone	· · · · · · · · · · · · · · · · · · ·	- 6	_					
Carbon Disulfide		- i	U					
1.1-Dichloroethene		107	×					
1.1-Dichloroethane		1	U	•				
1,2-Dichloroethene (t	otal)	1	U					
Chloroform		1	U					
1.2-Dichloroethane		1	U					
Z-Butanone		2	บ					
1,1,1-Trichloroethane		_ 1	U					
Carbon Tetrachloride_		1	U					
Vinyl Acetate		2	U					
Bromodichloromethane		_ j	Ü					
1,2-Dichloropropane_]	U					
cis-1.3-Dichloroprope	ле	- 1 - 91	Ü					
Trichloroethene Dibromochloromethane		- 1	X U					
		- ;	U					
1,1,2-Trichloroethane		- 691	· %					
Benzene Trans-1,3-Dichloropro	2000	- 23,	. X					
Bromoform	hene	- 1	Ŭ					
4-Hethy1-2-pentanone		- 2	Ü				•	
2-Hexanone		- 2	ប					
Tetrachloroethene		- 1	Ü					
1.1.2.2-Tetrachloroet	hane	- i	Ü					
*= Outside of EPA CLP		- 1	•					
- outside of FIX OF	40 1141634							

Xylene (total) ** Outside of EPA CLP QC limits.	Styrene	Ethylbenzene	Toluene
		9	4

RFW1: 926YEQ24-MB1

: 27:II : #6-II-L :

ZEZI BX:E230

Roy F. Weston, Inc. - Gulf Coast Laboratories
VOLATILES BY GC/MS. HSL LIST

RFW Batch Number:	92026419	Client:	Ek	VOLATILES BY G co <mark>Houseware</mark>	C/MS, HSL L		rk Order: 2994	Report Date: -02-02-0000	02/25/92 14:49 Page: 41
	Cust 10:	R-	5	VBLK	YBLK BS		VBLK	VBLK BS	VBLK
Sample	RFW#:	10	0	92GYE023-NB1	92GVE023-	HB1	92GVC040-HB1	92 GV C040-NB1	92GYE024-N81
Information	Matrix:	WATER		WATER	WATER		WATER	WATER	WATER
	D.f.:	1.	00	1.00	1.0		1.00	1.00	1.00
	Units:	ug/	L	ug/L	ug/l	L	ug/L	ug/L	ug/L
	To luene-d8	104	*	97 %	101	×	100 %	99 %	103 %
Surrogate Bro	mofluorobenzene	108	%	98 %	104	X.	102 %	107 %	104 %
Recovery 1,2-Di	chloroethane-d4	109	% f1	97 %	108	%	101 %	116 * %	94 %
Chloromethane		2	U	2 U	2	U	2 U	2 U	2 U
Bromomethane		2	U	2 U	2	U	2 U	2 U	2 U
Vinyl Chloride		9		1 U	ļ	U	1 U	1 U	1 U
Chloroethane		2	บ	2 U	2	U	2 U	2 U	2 U
Methylene Chloride		2		1 U	1	V	2	2 B	1 U
Acetone		Z	U	2 U	2	ľ	2	5 B	2 U
Carbon Disulfide		1	U	1 U	1	Ñ	1 U	1 U	1 1
1,1-Dichloroethene		Ī	U	1 U	114	X	1 0	132 %	iñ
1,1-Dichloroethane 1,2-Dichloroethene	7828211	1		1 U	ļ	Ü	ı u	1 U	1 U
1,2-vichiorbethene Chloroform	(1011)	12	11	1 0	į	U	1 U	1 U 1 U	1 0
1,2-Dichloroethane		1	IJ	1 0		U	1 U 1 U	i U	1 0
2-Butanone	······································	1	11	2 U	1	U U	2 U	1 U	2 11
1,1,1-Trichloroeth	204	1	11	1 0	2	11	ו ו	ו נ	1 11
Carbon Tetrachloric		i	ii	iŭ	i	II	iŭ	iŭ	iŭ
Vinyl Acetate	~ · · · · · · · · · · · · · · · · · · ·	;	ŭ	2 Ŭ	ż	il	2 Ŭ	2 0	2 11
Bromodichloromethan	ne .	7	ĬI.	ĩũ	i i	ŭ	īŭ	īū	i ü
1.2-Dichloropropane		i	ŭ	iŭ	i	ij	iŭ	iŭ	iŭ
cis-1,3-Dichloropro	opene	i	Ĭ	iŭ	i	1)	iũ	iū	iŭ
Trichloroethene		6		iŭ	97	į	iŭ	98 %	ĬŬ
Dibromochloromethan	ne	ĭ	U	iŭ	-1	Ũ	iŭ	ı ü	i v
1,1,2-Trichloroeth		i	ŭ	iŭ	ī	ŭ	iŭ	iŭ	ĨÙ
Benzene		ī	Ü	iŭ	80	Ť	iŭ	90 %	ĩũ
Trans-1,3-Dichloro	propene	3	Ū	i ŭ	1	ũ	iŭ	1 Ü	iŭ
Bronoform		i	Ū	iŭ	i	Ŭ	ĭŬ	i Ū	i Ū
4-Hethyl-2-pentanor	ne	2	ŭ	Ž Ũ	2	Ū	2 Ū	Ž Ū	ŽŪ
2-Hexanone	· · · · · · · · · · · · · · · · · · ·	2	Ū	ŽŪ	ž	Ū	ŽÜ	2 Ŭ	Ž Ū
Tetrachloroethene		1	ŭ	ĩũ	ĩ	Ŭ	ĭŬ	ĩ ǔ	ĩŬ
1,1,2,2-Tetrachlord	pethane	i	Ū	ìυ	i	Ŭ	ΙŪ	ĬŬ	īŪ
*= Outside of EPA (CIP OC Timits.	-	•		-	_	• •		- •

A STATE OF THE PARTY STATE OF TH	Cust 10:	7-1en 27-2KG R-5	VBLX	SB X78A	ABTX ABTX	SB XTBA	ABLK
	RFW#:	010	926VE023-MB1	926VE023-NB1 92GVE023-NB1	926YC040-MB1	926YC040-HB1 926YC040-HB1 926YE024-HB	92GVE024-MB1
Toluene		1 0	1	85 %	1 0	88	1 0
Chlorobenzene		_	1 U	96		100	
Ethylbenzene		1 6	1 U			. <u>_</u>	 : C
Styrene			_	_	· •	·	
<pre>xylene (total) *= Outside of EPA CLP QC limits.</pre>	QC limits.	_		-	-	-	-

SENT BY: E530

Roy F. Weston, Inc. - Gulf Coast Laboratories

DEM Dadah Musham. 09090410	VO	LATILES BY GC/N	r Coast Laborat 45, HSL LIST Hork	Order: 2994-0	Report Date: 0	2/25/92 14:49 Page: 3a
RFW Batch Number: 92026419	Client: Ekco		<u> </u>			
Cust ID:	L-3	L-4	L-4	L-4	L-4	L-5
Sample RFV#:	007	800	008 DL	008 MS	OO8 HSD	009
Information Matrix:	WATER	WATER	WATER	WATER	WATER	WATER
D.F.:	1.00	1.00	5.00	1.00	1.00	1.00
Units:	ug/L	ug/L	ug/L	ug/L	ug/L	ug/L
Taluene-d8	101 %	101 %	98 %	100 %	101 %	101 %
Surrogate Bromofluorobenzene	-104 %	107 %	108 %	106 %	107 %	106 %
Recovery 1,2-Dichloroethane-d4	97 %	126 * %	100 %	125 * %	127 * % f]=	118 * %
Chloromethane	2 U	S N	NA	2 U	2 Ú	2 U
Bromomethane	2 U	2 U	NA	2 U	2 U	2 U
Vinyl Chloride	1 U	ט נ	NA	1 U	1 U	37
Chloroethane	2 U	2 U	NA	2 U	2 U	2 U
lethylene Chloride	2 U	4 B U	NA	5 B	4 B	3 B
Acetone	2 U	2 U	NA	. 2 U	7 B	S N
Carbon Disulfide	1 U	1 U	KA	1 U	1 U	1 U 1 U
1.1-Dichloroethene	1 U	2 1	NA	113 %	120 %	?" 10 '
1.1-Dichloroethane	1 U	21 '	NA	21 _	21	
1,2-Dichloroethene (total)	1 U	E	20 '	E	t .	₹18
Chloroform	1 U	1 U	NA	1 U	1 V	1 0
1,2-Dichloroethane	1 U	1 บ	NA	1 U	1 U	1 U 2 U
2-Butanone	2 U	2 U	NA	2 U	2 U	• •
1,1,1-Trichloroethane	1 ย	1 U	NA	1 U	1 U	1 U 1 U
Carbon Tetrachloride	1 U	1 U	NA	1 U] U 2 U	2 0
Vinyl Acetate	2 U 1	2 U	MA	2 U	• •	1 0
Bromodichloromethane	1 U	1 U	KA	1 U	i n	1 0
1,2-Dichloropropane	1 U	1 U	NA	1 U	1 0	1 4
cis-1,3-Dichloropropene	1 U	1 U	NA.	1 U	1 0	1 0
Trichioroethene	1 U	1 U	NA	101 %	102 % 1 11	1 0
Dibromochloromethane	1 U	1 U	NA	i ñ	1 0	1 0
1,1,2-Trichlorgethane	1 U	1 U	NA	1 U		1 0
Benzene	1 U	1 U	NA	92 % 1 U	94 % 1 U	iu
Trans-1,3-Dichloropropene	1 A	ו נ	NA		1 0	iü
Bromoform	1 U	1 0	NA		1 U 2 U	2 U
4-Methy1-2-pentanone	2 U	2 U	NA	2 U	2 U	2 U
2-Hexanone	2 U	2 U	NA	2 U		
Tetrachloroethene	1 U	1 U	NA	j Ü	1 U	
1,1,2,2-Tetrachloroethane	1 U	1 U	NA	I U	1 U	1 U
- Outside of EPA CLP QC limits.						

Chlorobenzene Chlorobenzene Ethylbenzene Styrene Xylene (total) Xylene of E		
uene orobenzene ylbenzene ylbenzene ene (total) ene (total) Outside of EPA CLP QC limits.	RFU1:	925.019 Cust ID:
	007	iem kc
	800	-3 Ekc mansani
3333	10 800	Vor.
108 101 102 × ×	008 HS	North Wide - 2394-02-02-000
CCC**	OSK #500	1-4-10-10-10-10-10-10-10-10-10-10-10-10-10-
	909	

: 3**5**:11 : **56-**11-4 :

SENT BY: E580

RFW Batch Number: 921	026419	Client:	Ekco	Houseware	<u> </u>			Order: 2	994-0	leport Dai 2-02-0000		Page:	
	Cust ID:	Į-	1	L-1 Dug	•	L-1 Dug	P	L-1 Du	P	L-a	2	L-	.2
Sample	RFW#:	004 0	L	00!	5	005 DE	L	005 DI	L	000	5	006 E)L
Information	Matrix:	WATER		WATER		WATER		WATER		WATER		WATER	t
	D.F.:	40	.0	1.0	00	10.	.0	50	.0	1.0	00	20	0.0
	Units:	ug/	L	ug/i	•	ug/l	L	ug/l	L	ug/l	•	ug/	'L
	YoTuene-d8	102	*	99	×	100	*	99	*	103	*	99	7
Surrogate Bromof	luorobenzene	109	%	116	~	100	%	107	%	103	*	106	×
lecovery 1,2-Dichi	oroethane-d4	111	% f1-	130	* % :={}=:	97	% f3-	97 	% £1	101	% £1	99	%
hloromethane		NA	, , -	2	U	NA	, , -	NA	,,	2	Ü	NA	j
romomethane		NA		2	U	NA		NA		2	ีย	KA	
inyl Chloride		NA		23	•	NA		NA		1	U	NA	
hloroethane		NA		2	U	na		ŅA		2	U	NA	
ethylene Chloride_		NA		3	BU	NA		NA		2	u	KA	
cetone	·	NA		2	ีย	NA		NA		6	u	NA	
arbon Disulfide		NA		1	U	NA		NA		1	U	KA	
,1-Dichloroethene		NA		25		NA		NA		. 1	U	NA	
,1-Dichloroethane		NA	•		E	17		NA		12:	• • •	NA	
,2-Dichloroethene (t	iotal)	NA			E	110	•	NA		13		NA	
hloroform		NA		1	U	NA		NA		1	U	NA	
.2-Dichloroethane		, NA		1	U	NA		NA		1	Ų	NA	
-Butanone		NA		2	ຸປ	MA		MA		2	U	NA	
,1,1-Trichloroethane		NA		43		NA		NA		"5	•	NA	
arbon Tetrachloride		NA		1	Ü	NA		NA		1	ñ	NA	
Inyl Acetate		NA		Z	Ü	NA		NA		Z	U	NA	
romodichioromethane		NA NA		. 1	U	na Na		NA		1	t)	na Na	
.2-Dichloropropane		NA NA		1	Ü	NA NA		na Na		1	ŭ	NA NA	
is-1,3-Dichloroprope	:ne	'360		ı	E	ran.	E	1.430			E	710	
richloroethene ibromochloromethane				•	Ü	NA	Ε.			1	บ	NA NA	, ,
		na Na			Ŭ	NA NA		na Na			Ü	NA NA	
,1,2-Trichleroethane		NA NA		<u> </u>	บ	na NA		NA NA		1	ŭ	NA NA	
enzene	0000			1	Ŭ	KA		NA NA		1	ŭ	NA NA	
rans-1,3-Dichloropro	mene	MA		1	บ	na NA		NA NA		1	Ü	NA	
romoform_		NA		ı	บ					i	Ŭ	NA NA	*
-Methyl-2-pentanone		NA		2	U	MA		NA NA		2	U	na Na	
-Hexanone		NA		2,	_	AK		NA NA			ß		
etrachloroethene_		NA		1	Ŋ	MA		NA			U	NA NA	
,1,2,2-Tetrachloroet - Outside of EPA CLP		NA		1	U	NA		KA		1	V	NA	

WESTON- ROLE. RESION. LNC 1712-19

Toluene Chlorobe Ethylber Styrene Xylene Xylene			
Toluene Chlorobenzene Ethylbenzene Styrene Kylene (total) Kylene (total) Outside of EPA CLP QC limits.		Batter	
Eby Ci			
हैं।		CU 19	8
	RFW#:	Cust ID:	
3333	004 DL	T less	
	2		
	_	dng [-]	1
CCCC	005	- Cur	
****	005 DI	L-1 Pu	
	2	Work S	
ZZZZZ	005 DT	1-1 bup	
	7	994-02-	
900 900 900 900	96	DZ-0000	
cccc	√		
****	006 DT	Page:	1
-		3	



WESTON-GULF COAST LABORATORIES. NC.

2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 | 219) 385-7077 | 815) 723-752

ANALYT DAL ERRORT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 5, 1989

RE: Grdwater L-1-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88 GCL#: 144693

VOLATILE COMPOUNDS	r e sul t	DETECTION LIMIT
1,2-Dichloropropane	BDL	5 ug/1 U
Cis-1,3-Dichloropropene	BDL	5 ug/l U
Trichloroethene	-210 ·	5 ug/l (e')
Dibromochloromethane	BDL	5 ug/l U
1,1,2-Trichloroethane	BDL	5 ug/1 U
Benzene	BDL	5 ug/1 U
Trans-1,3-Dichloropropene	BDL	5 ug/1 U
Bromoform	BDL	5 ug/1 U
4-Methyl-2-Pentanone	BDL	10 ug/l U
2-Bexanone	BDL	10 ug/l U
Tetrachloroethene	BDL	5 ug/l U
1,1,2,2-Tetrachloroethane	BDL	5 ug/l 0
Toluene	BDL	5 ug/l U
Chlorobenzene	BDL	5 ug/l U
Ethylbenzene	BDL	5 ug/l U
Styrene	BDL	5 ug/l U
Xylene	BDL	5 ug/1 U
•		



WESTON-GULF COAST LABORATORIES. NC 2417 Band St., University Park, Illinois 60466

Phones: -3121 534-5290 | 2191 \$85-7977 | 815-123-13

4441 7 041 482047

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 5, 1989

RE: Grdwater L-1-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88

Date Received: 11/18/88

GCL#:

144693

VOLATILE COMPOUNDS	RESULT	DETEC	TION I	LIMIT
Chloromethane	BDL	10	ug/1	Ū
Bromomethane	BDL	10	ug/l	Ū
Vinyl Chloride	48 ,	10	ug/l	······································
Chloroethane	BDL.	10	ug/l	Ū
Methylene Chloride	30	5	ug/1	B
Acetone	11	10	ug/l	B
Carbon Disulfide	BDL	5	ug/1	Ū
1,1-Dichloroethene	3	5	ug/1	J
1,1-Dichloroethane	:57	5	ug/l	
1,2-Dichloroethene (total)	61	5	ug/l	··
Chloroform	BDL	5	ug/1	Ū
1,2-Dichloroethane	BDL	5	ug/1	· 0
2-Butanone	BDL	10	ug/1	0
1,1,1-Trichloroethane	49	5	ug/1	
Carbon Tetrachloride	BDL	5	ug/1	Ū
Vinyl Acetate	BDL	10	ug/l	O
Bromodichloromethane	BDL	5	ug/l	0



WESTON-GULF COAST LABORATORIES, INC. 2417 Bong St., University Park, illinois 60466

Phones: 312) 534-5200 | 219) 385-7077 | 816) 700-153:

AV4_/T [4] PEPCAT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater I-5-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/17/88

Date	Received:	11/18/88
GCL#	:	144692

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U



WESTON-GULF COAST LABORATORIES INC. 2417 Bond St., University Park, Hinois 50466

Phones: .312) 534-5290 | 219: 385-7977 | 815) 123-150

ANALYTICAL ABACAT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 5, 1989

RE: Grdwater I-5-GW-001

American Home Products

W.O. \$ 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88

GCL#: 144692

VOLATILE COMPOUNDS	result	DETECT	!ION	LIMIT
1,2-Dichloropropane	BDL	5	ug/l	U
Cis-1,3-Dichloropropene	BDL	5	ug/l	L Ū
Trichloroethene	₹40	5	ug/1	e
Dibromochloromethane	BDL	5	ug/l	נ ס
1,1,2-Trichloroethane	BDL	5	ug/l	L U
Benzene	6	5	ug/l	L J
Trans-1,3-Dichloropropene	BDL	5	ug/l	U
Bromoform	BDL	5	ug/l	L Ū
4-Methyl-2-Pentanone	BDL	10	ug/l	l U
2-Hexanone	BDL	10	ug/l	L U
[etrachloroethene	BDL	5	ug/]	1 0
1,1,2,2-Tetrachloroethane	BDL	5	ug/l	L 0
Toluene	BDL	5	ug/1	U
Chlorobenzene	BDL	5	ug/l	U
Ethylbenzene	BDL	5	ug/l	l U
Styrene	BDL	5	ug/]	L O
Xylene	BDL	5	ug/	1 0



WESTON-GULF COAST LABORATORIES, INC 2417 Bond St., University Park, Illinois 50466

Phones: 3121 534-5200 (219) 385-7077 (815) 720-7500

ANALYT DAU REROAT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 5, 1989

RE: Grdwater I-5-GW-001
American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88
GCL#: 144692

VOLATILE COMPOUNDS	RESULT	DETECTI	ON LIHIT
Chloromethane	BDL	10 u	g/l U
Bromomethane	BDL	10 u	g/l U
Vinyl Chloride	BDL	10 u	g/l U
Chloroethane	BDL	10 Ա	g/1 U
Methylene Chloride	31	5 u	g/1 B
Acetone	26	10 u	g/1 B
Carbon Disulfide	BDL	5 u	g/l U
1,1-Dichloroethene	5	5 u	g/l
1,1-Dichloroethane	~B6	5 u	g/l
1.2-Dichloroethene (total)	BDL	5 u	g/l U
Chloroform	BDL	5 u	g/l U
1,2-Dichloroethane	BDL	5 u	g/1 0
2-Butanone	BDL	10 u	g/1 U
1,1,1-Trichloroethane	€40	5 u	g/1
Carbon Tetrachloride	BDL	5 u	g/1 U
Vinyl Acetate	BDL	10 u	g/l U
Bromodichloromethane	BDL	5 u	g/l U
	·		



WESTON-GULF COAST LABORATORIES INC. 241T Bond St., University Park, Illinois 30466 Phones: (312) 534-5200 (219) 385-707T | 315-723-758

LNALYT GAL REPORT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater I-2-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/17/88

Date Received: 11/18/88

GCL#:

144691DL

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT	
1,2-Dichloroethene (total)	3880	50 ug/l D	
Trichloroethene	-1 0	50 ug/l D	
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WESTON-GULF COAST LABORATORIES, INC. 2417 Bond St., University Park, Illinois 50466 Phones: 312) 534-5290 (213) 885-7077 (815) 723-752

7/77\1.077 BEb0et

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater I-2-GW-001

American Home Products W.O.# 2994-02-03

Sample Date: 11/17/88

Date Received: 11/18/88 GCL#: 144691

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U
		·



WESTON-GULF COAST LABORATORIES. NC. 2417 Bond St., University Park, Illinois 60486

Phones: (3*2) 534-5200 (219) 885-707** (815) 723-753

אלק"ים נין שבשנטי

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater I-2-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88

GCL#: 144691

VOLATILE COMPOUNDS	result	DETEC	TION I	IMIT
1,2-Dichloropropane	BDL	5	ug/l	σ
Cis-1,3-Dichloropropene	BDL	5	ug/l	Ū
Trichloroethene	680	5	ug/l	©
Dibromochloromethane	BDL	5	ug/l	Ū
1,1,2-Trichloroethane	BDL	5	ug/1	U
Benzene	BDL	5	ug/l	Ū
Trans-1,3-Dichloropropene	BDL	5	ug/l	σ
Bromoform	BDL	5	ug/l	ט
4-Methyl-2-Pentanone	BDL	10	ug/l	Ū
2-Bexanone	BDL	10	ug/l	Ū
Tetrachloroethene	BDL	5	ug/l	σ
1,1,2,2-Tetrachloroethane	BDL	5	ug/1	U
Toluene	BDL	5	ug/l	Ū
Chlorobenzene	BDL	5	ug/l	U
Ethylbenzene	BDL	5	ug/l	Ū
Styrene	BDL	5	ug/l	ט
Xylene	BDL	5	ug/l	Ū
				



WESTON-GULF COAST LABORATORIES NO. 2417 Bond St., University Park, Illinois 50466

Phones: 312) 534-5200 (213) 385-7077 (815) 123-753:

4 V 4 L Y T C 4 L 4 E 4 C A 7

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater I-2-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88 GCL#: 144691

VOLATILE COMPOUNDS	RESULT	DETEC	TION	LIHIT
Chloromethane	BDL	10	ug/l	Ū
Bromomethane	BDL	10	ug/l	ס
Vinyl Chloride	56	10	ug/l	
Chloroethane	BDL	10	ug/l	U
Methylene Chloride	34	5	u g /1	B
Acetone	34	10	ug/l	B
Carbon Disulfide	BDL	5	ug/l	Ū
1,1-Dichloroethene	#22	5	ug/l	
1,1-Dichloroethane	£ <u>£</u> 0	5	ug/l	
1,2-Dichloroethene (total)	300	5	ug/l	•
Chloroform	BDL	5	ug/l	Ū
1,2-Dichloroethane	BDL	5	ug/l	Ū
2-Butanone	BDL	10	ug/1	Ū
1,1,1-Trichloroethane	57	5	ug/l	
Carbon Tetrachloride	BDL	5	ug/l	σ
Vinyl Acetate	BDL	10	ug/1	Ū
Bromodichloromethane	BDL	5	ug/l	U



#55TOM-GULFICIAST_LABORATOR(ES_NC)
241T Bond 3: University Park, Illinois 30466
Phones: 012 304-3200 | 0131-335-707T | 8 (5 100173)

a wi Till agains

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater D-4-30-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/17/88

Date Received: 11/18/88

GCL#:

144690DL

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
1,1,1-Trichloroethane	\$200 0	10000 ug/l D
Trichloroethene	9220000	10000 ug/l D
		
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WESTON-GULF DO AST LABORATORIES INC 1417 Bond St. University Park (Ondis \$0486 Phones: 072) \$04-\$200 | 219 \$45-7077 | 318 720-720

42 7 143 823115

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater D-4-30-GW-001 American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88

GCL#:

144690

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U
	- 	



WESTON-GULF TOAST LABORATORIES NO. 1417 Bond St. University Park Land's \$0456. Phones 012: \$24 \$200 | 219 \$450,7077 | 3150,720 15

- ----

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater D-4-30-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88

GCL#: 144690

VOLATILE COMPOUNDS	RESULT	DETEC	TION I	LIMIT	
1,2-Dichloropropane	BDL	5	ug/l	Ū	
Cis-1,3-Dichloropropene	BDL	5	ug/l	Ū	
Trichloroethene	1900	5	ug/l	e	
Dibromochloromethane	BDL	5	ug/1	Ū	
1,1,2-Trichloroethane	53.40	5	18/ 1		
Benzene	BDL	5	1g/l	ี่	
Trans-1,3-Dichloropropene	BDL	5	ug/1	J	
Broacform	BDL	5	ug/l	1	
4-Methyl-2-Fentanone	: T.	10	1g/l	J	
2-Bexanone	EDL	10	1g,/1	IJ	
Tetrachloroethene	.75.5	5	ug/1		
1,1,2.2-Tetrachloroethane	BDL	5	1g/l	ס	
Toluene	130	5	ug/1		
Chlorobenzene	BDC	5	ug/l	ס	
Ethylbenzene	3	5	ug/1	J	
Styrene	BDL	5	ug/1	Ū	
Xylene (total)	i3 :	5	ug/1		



WESTON-GULF COAST LABORATORIES INC 2417 Band St., University Parx, Minous \$0466 Phones: 3121 534-5200 | 2191 385: 7077 | 315: 703-752

A .ALYT [AL RED] AT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater D-4-30-GW-001 American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88
GCL#: 144690

VOLATILE COMPOUNDS	RESULT	DETECT	TION LIMIT
Chloromethane	BDL	10	ug/l U
Bromomethane	BDL	10	ug/l U
Vinyl Chloride	(3)	10	ug/l J
Chloroethane	BDL	10	ug/l U
Methylene Chloride	59	5	ug/1 B
Acetone	82	10	ug/l B
Carbon Disulfide	BDL	5	ug/l U
1.1-Dichloroethese	1900	5	ug/l e
1,1-Dichloroethane	1,000	5	ug/l (e)
1.2-Dichloroethene (total)	270	5	1g/l e
Chloroform	10	5	ug/l
1,2-Dichloroethane	₹3	5	ug/l
2-Butarone	BDL	10	ug/l U
1,1,1-Trichloroethane	3000	5	ug/l (e)
Carbon Tetrachloride	BDL	5	ug/l U
Vinyl Acetate	BDL	10	ug/l 0
Bromodichloromethane	BDL	5	ug/l U



₩ESTON-GULF COAST LABORATOR ET INC. 2417 Eand St., University Park (1997) (1997)

Phones: (312) 534-5200 (219) 885-7077 (815 723 75)

ANALYT DAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 4, 1989

RE: Ground Water L-5-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/16/88
Date Received: 11/17/88

GCL#: 144615

VOLATILE COMPOUNDS	RESULT	DETEC	TION [IMIT
2-Chloroethyl vinyl ether	BDL	10	ug/l	Ū
1,3-Dichlorobenzene	BDL	10	ug/l	Ū
1,2-Dichlorobenzene	BDL	10	ug/l	Ū
1,4-Dichlorobenzene	BDL	10	ug/l	U
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WESTON-GULF COAST LABORATORIES INC. CATT BORD St. Long. 1. To a control 64464

Phones: (312) 534-5200 (219) 885-7077 (815-723 TFT

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 4, 1989

RE: Ground Water L-5-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/16/88 Date Received: 11/17/88

GCL#:

144615

VOLATILE COMPOUNDS	RESULT	DETECT	'ION L	IMIT
1,2-Dichloropropane	BDL	5	ug/l	U
Cis-1,3-Dichloropropene	BDL	5	ug/l	Ū
Trichloroethene	BDL	5	ug/l	Ū
Dibromochloromethane	BDL	5	ug/l	Ū
1,1,2-Trichloroethane	BDL	5	ug/l	Ü
Benzene	BDL	5	ug/l	Ū
Trans-1,3-Dichloropropene	BDL	5	ug/l	Ū
Bromoform	BDL	5	ug/l	U
4-Methyl-2-Pentanone	BDL	10	ug/l	Ū
2-Hexanone	BDL	10	ug/l	Ū
Tetrachloroethene	BDL	5	ug/l	Ū
1,1,2,2-Tetrachloroethane	BDL	5	ug/l	Ü
Toluene	BDL	5	ug/l	Ū
Chlorobenzene	BDL	5	ug/l	Ū
Kthylbenzene	BDL	5	ug/l	Ū
Styrene	BDL	5	ug/l	Ū
Xylene	BDL	5	ug/l	Ū



WESTON-GULF COAST LABORATORIES INC. 2417 Bond St. Utilis 1994 444

Phones: (312) 534-5200 (219) 885-7077 (815, 723-752)

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 4, 1989

RE: Ground Water L-5-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/16/88 Date Received: 11/17/88

VOLATILE COMPOUNDS	RESULT	DETEC	TION LIMIT
Chloromethane	BDL	10	ug/l U
Bromomethane	BDL	10	ug/l U
Vinyl Chloride	510	10	ug/l
Chloroethane	BDL	10	ug/l U
Methylene Chloride	62	5	ug/1 B
Acetone	74	10	ug/l B
Carbon Disulfide	BDL	5	ug/l U
1,1-Dichloroethene	BDL	5	ug/l U
1,1-Dichloroethane	7-8	5	ug/l J
1,2-Dichloroethene (total)	-9 2	5	ug/l
Chloroform	BDL	5	ug/l U
1,2-Dichloroethane	BDL	5	ug/l U
2-Butanone	BDL	10	ug/l U
1,1,1-Trichloroethane	BDL	5	ug/l U
Carbon Tetrachloride	BDL	5	ug/l U
Vinyl Acetate	BDL	10	ug/l U
Bromodichloromethane	HDL	5	ug/l U



2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-7533

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

- I SINGLAN

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-4-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/16/88 Date Received: 11/17/88 GCL#: 144614

VOLATILE COMPOUNDS	result	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	1.00 ug/l U
1,3-Dichlorobenzene	BDL	0.08 ug/l U
1,2-Dichlorobenzene	BDL	0.09 ug/l U
1,4-Dichlorobenzene	BDL	0.09 ug/l U
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2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-75.

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-4-GW-001

American Bome Products

W.O.# 2994-02-03

Sample Date: 11/16/88
Date Received: 11/17/88
GCL#: 144614

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
1,2-Dichloropropane	BDL	0.24 ug/l U
Cis-1,3-Dichloropropene	BDL	0.1 ug/l U
Trichloroethene	BDL	0.27 ug/l U
Dibromochloromethane	BDL	0.18 ug/l U
1,1,2-Trichloroethane	BDL	0.20 ug/l U
Benzene	BDL	0.07 ug/l U
Trans-1,3-Dichloropropene	BDL	0.10 ug/l U
Bromoform	BDL	0.15 ug/l U
4-Methyl-2-Pentanone	BDL	1.0 ug/l U
2-Bexanone	BDL	1.0 ug/l U
Tetrachloroethene	BDL	0.19 ug/l U
1,1,2,2-Tetrachloroethane	BDL	0.23 ug/l U
Toluene	0.25	0.30 ug/l J
Chlorobenzene	BDL	0.18 ug/1 U market
Sthylbenzene	BDL	0,19 mg/1 - 0
Styrene	BDL	0.18 ug/1 0
lylene	BDL	0.10 w/. 3



2417 Bond St., University Park, Linois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-753

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

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ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-4-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/16/88
Date Received: 11/17/88
GCL#: 144614

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
Chloromethane	BDL	0.41 ug/l U
Bromomethane	BDL	0.27 ug/l U
Vinyl Chloride	BDL	0.23 ug/l U
Chloroethane	BDL	0.14 ug/l U
Methylene Chloride	1.3	0.5 ug/l B
Acetone	4.4	1.0 ug/l
Carbon Disulfide	4 .1	0.5 ug/l
1,1-Dichloroethene	BDL	0.26 ug/l U
1,1-Dichloroethane	°50	0.17 ug/l
1,2-Dichloroethene (total)	BDL	0.17 ug/l U
Chloroform	BDL	0.14 ug/l U
1,2-Dichloroethane	BDL	0.1 ug/l U
2-Butanone	BDL	1.0 ug/l U
1,1,1-Trichloroethane	BDL	0.5 ug/1 U
Carbon Tetrachloride	EDL	0.08 ug/1 U
Vinyl Acetate	BDL	1.0 ug/l U
Bromodicatorometophe	BDL	0.52 ug/1 U.



2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-753.

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample S-7-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144532DL

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VOLATILE	COMPOUNDS	RESU	LT DETI	CTION	LIMIT
1,1-Dichl	oroethane	43	5	ug/l	D
1,1,1-Tri	chloroethane	.156	5	ug/l	D
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2417 Bond St., University Park, Illinois 50466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-752

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample S-7-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144532

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VOLATILE COMPOUNDS	result	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	1.00 ug/l U
1,3-Dichlorobenzene	BDL	0.08 ug/l U
1,2-Dichlorobenzene	BDL	0.09 ug/l U
1,4-Dichlorobenzene	BDL	0.09 ug/l U

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2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-753

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample S-7-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/14/88
Date Received: 11/16/88
GCL#: 144532

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VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
1,2-Dichloropropane	BDL	0.24 ug/l U
Cis-1,3-Dichloropropene	BDL	0.1 ug/l U
Trichloroethene	46	0.27 ug/l
Dibromochloromethane	BDL	0.18 ug/l U
1,1,2-Trichloroethane	BDL	0.20 ug/l U
Benzene	47 15	0.07 ug/l
Trans-1,3-Dichloropropene	BDL	0.10 ug/l U
Bromoform	BDL	0.15 ug/l U
4-Methyl-2-Pentanone	BDL	1.0 ug/l U
2-Hexanone	BDL	1.0 ug/l U
Tetrachloroethene	BDL	0.19 ug/l U
1,1,2,2-Tetrachloroethane	BDL	0.23 ug/l U
Toluene	13	0.30 ug/l
Chlorobenzene	BDL .	0.18 ug/1 0
Ethylbenzene	0.29	0:19 ug/1
Styrene	BDL .	0.18 us/1 U
Xylene		Carle Water



2417 Bond St., University Para minois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-753

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample S-7-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144532

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VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
Chloromethane	BDL	0.41 ug/l U
Bromomethane	BDL	0.27 ug/l U
Vinyl Chloride	BDL	0.23 ug/l U
Chloroethane	BDL	0.14 ug/l U
Methylene Chloride	0.79	0.5 ug/l B
Acetone	.17	1.0 ug/l
Carbon Disulfide	BDL	0.5 ug/l D
1,1-Dichloroethene	11.9	0.26 ug/l
1,1-Dichloroethane	X	0.17 ug/l (e)
1,2-Dichloroethene (total)	BDL	0.17 ug/l U
Chloroform	BDL	0.14 ug/l U
1,2-Dichloroethane	BDL	0.1 ug/l U
2-Butanone	BDL	1.0 ug/l U
1,1,1-Trichloroethane	100	0.5 ug/l 6B
Carbon Tetrachloride	BDL	0.08 ug/1 U
Vinyl Acetate	BDL	1:0 ug/1 U- *-
Bromodichloromethane	BDL	0.62 ug/1 U



2417 Bond St., University Park, Empis 50466

Phones: (312) 534-5200 (219) 885-7077 (815 723 7

ANALYT CAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample OWS4-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/15/88 Date Received: 11/16/88 GCL#: 144519

VOLATILE COMPOUNDS	RESULT	DETECTION LIM	T
2-Chloroethyl vinyl ether	BDL	1.00 ug/l (j
1,3-Dichlorobenzene	BDL	0.08 ug/l [)
1,2-Dichlorobenzene	BDL	0.09 ug/l (J
1,4-Dichlorobenzene	BDL	0.09 ug/l (J
			
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2417 Bond St., University Park, Illinois 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-7531

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample OWS4-GW-001

A PROPERTY OF THE PROPERTY OF

American Home Products

W.O.# 2994-02-03

Sample Date: 11/15/88
Date Received: 11/16/88
GCL#: 144519

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
1,2-Dichloropropane	BDL	0.24 ug/l U
Cis-1,3-Dichloropropene	BDL	0.1 ug/l U
Trichloroethene	0.16	0.27 ug/l J
Dibromochloromethane	BDL	0.18 ug/l U
1,1,2-Trichloroethane	BDL	0.20 ug/l U
Benzene	0.37	0.07 ug/l
Trans-1,3-Dichloropropene	BDL	0.10 ug/l U
Bromoform	BDL	0.15 ug/l U
4-Methyl-2-Pentanone	BDL	1.0 ug/l U
2-Bexanone	BDL	1.0 ug/l U
Tetrachloroethene	RDL	0.19 ug/l U
1,1,2,2-Tetrachloroethane	BDL	0.23 ug/l U
Toluene	0.80	0.30 ug/l
Chlorobenzene	BDL	0.18 ug/l I
Ethylbenzene	0.33	0.19 ug/l
Styrene	BDL	0.18 ug/1 0
Xylene	0.42	The state of the s



2417 Bond St., University Park, Hillinois 50466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-*53

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATIN: Ms. Sherrerd Steele

DATE: January 1989

RE: Ground Water Sample OWS4-GW-001

American Bome Products

W.O.# 2994-02-03

Sample Date: 11/15/88 Date Received: 11/16/88 GCL#: 144519

VOLATILE COMPOUNDS	result	DETECTION LIMIT	
Chloromethane	BDL	0.41 ug/l U	
Bromomethane	BDL	0.27 ug/l U	
Vinyl Chloride	3.5	0.23 ug/l	
Chloroethane	BDL	0.14 ug/l U	
Methylene Chloride	1.9	0.5 ug/l B	
Acetone	7.1	1.0 ug/l	
Carbon Disulfide	2.8	0.5 ug/l	
1,1-Dichloroethene	BDL	0.26 ug/l U	
1,1-Dichloroethane	BDL	0.17 ug/l U	
1,2-Dichloroethene (total)	BDL	0.17 ug/l U	
Chloroform	BDL	0.14 ug/l U	
1,2-Dichloroethane	BDL	0.1 ug/l U	
2-Butanone	BDL	1.0 ug/l U	
1,1,1-Trichloroethane	0.21	0.5 ug/l B J	
Carbon Telrachioride	BDL	0.08 ug/E-034	loi
Vinyl Acetate	BDL	1.0° ug/l U	
Bromod (ed (cross) date	A STATE OF THE STA	10.52 uzy uz	-



2417 Bond St., University Park Himbis 60466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-7533

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-7-GW-001

American Bome Products

W.O.# 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144516DL

VOLATILE COMPOUNDS	RESULT	DETECTI	ON LIM	IT
1,1-Dichloroethene	160	5	ug/l	D
1,1-Dichloroethane	₹200	500	ug/l	D
1,1,1-Trichloroethane	440	500	ug/l	D
Trichloroethene	830	500	ug/l	D
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1417 Bond St., University Park, Illinois 60456

Phones: (312) 534-5200 (219) 885-7077 (815) 723 7532

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-7-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144516

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	1.00 ug/l U
1,3-Dichlorobenzene	BDL	0.08 ug/l U
1,2-Dichlorobenzene	BDL	0.09 ug/l U
1,4-Dichlorobenzene	BDL	0.09 ug/l U
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2417 Sond St., University Park, Illinois 50456

Phones: (312) 534-5200 (219) 885-7077 (815) 723-7533

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RK: Ground Water Sample I-7-GW-001

American Home Products

W.O.# 2994-02-03

The state of the state of

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144516

RESULT	DETECTION LIMIT
BDL	0.24 ug/l U
BDL	0.1 ug/l U
110	0.27 ug/l 📵
BDL	0.18 ug/l U
7.4 1	0.20 ug/l
BDL	0.07 ug/l U
BDL	0.10 ug/l U
BDL	0.15 ug/l U
BDL	1.0 ug/l U
BDL	1.0 ug/l U
- 0.88	0.19 ug/l
BDL	0.23 ug/1 U
1.9	0.30 ug/l
BDL	0.18 ug/l U
0.22	0.19 vg/1
BDL	0.18 ug/1 U
BDG	STATE OF THE STATE
	BDL BDL 7.4 BDL BDL BDL BDL BDL BDL BDL BD

Company Company (Company)



2417 Bond St. University Park, Juniors 50466

Phones: (312) 534-5200 (219) 885-7077 (815) 720 75

1988 DATA

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

All the second

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water Sample I-7-GW-001

American Bome Products

W.O.# 2994-02-03

Sample Date: 11/14/88 Date Received: 11/16/88 GCL#: 144516

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
Chloromethane	BDL	0.41 ug/l U
Bromomethane	BDL	0.27 ug/l U
Vinyl Chloride	5.3	0.23 ug/l
Chloroethane	6.4	0.14 ug/l
Methylene Chloride	1.7	0.5 ug/l B
Acetone	8.1	1.0 ug/l
Carbon Disulfide	BDL	0.5 ug/l U
1,1-Dichloroethene	120	0.26 ug/l
1,1-Dichloroethane	290	0.17 ug/l (e)
1,2-Dichloroethene (total)	16	0.17 ug/l
Chloroform	EDL.	0.14 ug/l U
1,2-Dichloroethane	3.7.	0.1 ug/l
2-Butanone	2.6	1.0 ug/l
1,1,1-Trichlorgethane	190	0,5 ug/1 (6.)
Carbon Tetrachloride.	BOL	0.08 ug/1- B
Vinyl Acetate	BDL	1.0 ug/l U
Bromodichlogogetaine	BOL	0.52 u/1 U

Chlorobenzene Ethylbenzene

45)	F. Westor The. Wall freest	entraterris -	
	VOLATILES BY GC/MS. HSL		Report D

Client: Ecko Housewares Hork Order: 2994-02-03-0035 Page: 4a RFW Batch Number: 91086851 VELK BS ABLK Cust ID: VBI K Sample RFW#: 918YB297-MB1 916VE300-KB1 918VR300-MR1 VATER Information WATER WATER Matrix: 1.00 1.00 1.00 D.F.: ug/L ug/L Units: uq/L To luene-dB 100 Bromofluorobenzene Ľ 103 96 101 Surrogate 1.2-Dichloroethane-d4 97 × 88 97 Recovery Chloromethane u u u 2 Bromomethane U Vinyl Chloride U Ch loroethane Methylene Chloride IJ Acetone Carbon Disulfide U 1.1-Dichloroethene 51 U 1,1-Dichloreethane 1,2-Dichiaroethene (total) Chloroform 1,2-Dichloroethane 2-Butanone U 1.1.1-Trichloroethane Carbon Tetrachloride_____ Vinyl Acetate Bromodichioromethane 1.2-Dichloropropene cis-1,3-Dichloropropene____ U Trichloroethene Dibromochioromethane 1.1.2-Trichloroethane ŧ, Trans-1,3-Dichloropropene U IJ Bromoform 4-Methyl-2-pentanone

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1,1,2,2-Tetrachloroethane *= Outside of EPA CLP QC limits.

2-Hexanone

Tetrach loroethene



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RFW Batch Number: 91086861	Client: Ecko	Housewares	Work	Order: 2994-02	-03-0035	Page: 3b K
Cust ID:	L-4	L-4	L-6	L-5	R-5	R-5
RFW#:	008 HS	008 MSD	009	009 DL	010	010 DL 👸
Toluene	75 * %	77 %	1 0	NA .		Į
Chlorobenzene	101 %	101 %	i ü	NA	i Ū	NA
Ethyl benzene	1 U	1 U	1 - U	NA	ĪŪ	NA
Styrene	1 U	1 U	1 U	NA	1 U	NA
Xylene (total)	1 U	1 U	1 U	NA	i v	NA

F. ASTONIANCE OF COME AND LEST PROPERTY DATE: 09/12/91 16:

FM Batch Number: 91	1000351	Client:	Ecko	Houseware	\$		lork D	rder: 29	<u>94-0</u> 2	-03-0035		Page:	3a
	Cust ID:	L-4	}	L-4	,	L- <u>5</u> (237)		L-5	i	R-5		R-5	•
ample	RFW#:	008 MS	:	OOS MSD		009		009 DL		010	,	010 DL	
nformation	Matrix:	WATER		WATER		WATER		WATER		WATER		WATER	
	D.F.:	1.0	00	1.0	0	1.00)	5.0	0	1.0	0	5.0	0
	Units:	ug/l	•	ug/L		ug/L		ug/L	•	ug/L		ug/l	-
	Toluene-d8	90	7	91	*	98	*	109	×	95	*	103	*
rrogate Bromo	fluorobeszene	105	%	104	*	95	%	92	*	101	*	95	*
covery 1,2-Dich	loroethane-d4	96	%	91	% -61	92	% .£1	94 	% f1	92	% _f]	103	% f1
loromethane		2	v	2	U	2	V	NA	,,	2	U	NA	,,
omomethane		2	U	2	Ü	2	U	NA		2	U	NA	
nyl Chloride		5		5			E 120	1150	•		Esi	51	
loroethane		2	V	2	ช	2	U	NA		2	Ų	NA	
thylene Chloride_		0.6	JB	0.B	JB		JBU	NA		1	B	NA	
:etone		2	U	2	V	2	U	NA		2	U	NA	
rbon Disulfide		1	U	1	U	1	V	NA		1	U	NA	
1-Dichloroethene		53 4	* %	54 ^	*	1	U	NA		1	U	NA	
1-Dichloroethane		19		21		শ8 '	•	NA		4		NA	
2-Dichloroethene	totall	64		67			E 67	163	•		ESY	54	
loroform		1	U	1	U	1	U	NA		1	U	NA	
2-Dichloroethane		i	Ũ	ì	Ü	Ī	V	NA		1	U	NA	
Butanone		ż	ŭ	Ž	Ŭ	Ž	Ũ	NA		Ž	Ū	NA	
1,1-Trichloroethan	16	ī	ũ	ī	Ŭ	ī	Ũ	NA		ī	Ū	NA	
rbon Tetrachloride	` <u> </u>	ī	Ü	ī	Ū	i	Ū	NA		ī	Ū	NA	
inyl Acetate		ż	Ŭ	Ž	Ū	2	Ü	NA		2	U	NA	
romodichloromethane	<u> </u>	ī	ŭ	ĭ	Ŭ	ĭ	Ŭ	NA		ī	Ü	NA	
2-Dichloropropane		i	ŭ	ī	Ũ	i	Ŭ	NA		Ĭ	Ŭ	NA	
s-1,3-D1chloroprop	iene	i	Ŭ	i	ŭ	i	ij	NA		ī	Ũ	NA	
richioroethene		84	%	86	Ÿ.	i	Ŭ	NA		17	_	NA	
bromochloromethane		07 1	ű	1	ũ	1	Ŭ	NA		'n	U	NA	
1,2-Trichloroethan	<u> </u>	1	Ü	1	Ŭ	1	Ŭ	NA		i	บั	NA	
	· · · · · · · · · · · · · · · · · · ·	76	%	75 4	_	1	Ü	NA NA		•	Ŭ	NA NA	
enzene rans-1,3-Dichloropr	Mana	10	76 U	79,"	W U	1	บ ม	NA NA		,	Ü	NA NA	
-ans-1,3-uicnioropr	ohene		u	i.	Ü	1	U	na NA		, , , , , , , , , , , , , , , , , , ,	Ŭ	NA NA	
romoform		Ĭ	_	i	fi O	l 2	11			1	IJ	NA NA	
-Methy1-2-pentanone		2	Ü	(U	ζ.	U H	NA		2	Ü	NA NA	
-Hexanone		2	Ü	2	-	7	_	NA		2	U U	NA NA	
etrachloroethene	10.	Ĭ	Ü	į	U	i	U	NA			_	•	
.1.2.2-Tetrachloroe	Inane		U	1	U	I	U	NA		1	U	NA	

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WESTON-GULF COAST LABORATORIES, INC. 241" Bond St., University Park, Hinois 60466

Phones: (312) 534-5200 (219) 385-7977 (815) 720-753

ANALYTICAL REPORT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 5, 1989

RE: Grdwater L-1-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88

GCL#:

144693

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U
		



WESTON-GULF COAST LABORATORIES. NC. 241" Bond St., University Park, Illinois 50466 Phones: (312) 534-5290 (219) 385-7077 (815) 720-752

AVALIT DAL REPORT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater L-2-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88

RESULT	DETEC	TION L	IMIT
BDL	10	ug/l	Ū
BDL	10	ug/l	U
BDL	10	ug/l	U
BDL	10	ug/l	ס
31	5	ug/ 1	<u>B</u>
10	10	u g /1	B
BDL	5	u g /1	Ū
BDL	5	ug/1	Ū
BDL	5	ug/1	σ
BDL	5	ug/l	Ū
BDL	5	ug/1	U
BDL	5	ug/l	Ū
BDL	10	ug/l	σ
*26	5	ug/l	
BDL	5	u g /l	U
BDL	10	ug/l	σ
BDL	5	ug/l	U
	BDL	BDL 10 BDL 10 BDL 10 BDL 10 BDL 10 31 5 10 10 BDL 5 BDL 10	BDL 10 ug/l BDL 10 ug/l BDL 10 ug/l BDL 10 ug/l 31 5 ug/l 10 10 ug/l BDL 5 ug/l BDL 10 ug/l BDL 10 ug/l BDL 10 ug/l BDL 5 ug/l BDL 10 ug/l



WESTON-GULF COAST LABORATORIES NC. 2417 Bond St., University Park, Minois 50466

Phones: 312) \$34-5200 | 219) 385-737" (815) 723-752.

ANALYT DAL REPORT

GCL#:

TO: Roy F. Weston, Incorporated

1 Weston Way

Ethylbenzene

Styrene

Xylene

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater L-2-GW-001
American Home Products

W.O.# 2994-02-03

5 ug/l 0

ug/1

ug/l

5

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Sample Date: 11/17/88
Date Received: 11/18/88

144694

VOLATILE COMPOUNDS	RESULT	DETEC	TION	LIMIT
1,2-Dichloropropane	BDL	5	ug/l	Ū
Cis-1,3-Dichloropropene	BDL	5	ug/l	U
Trichloroethene	130	5	ug/l	
Dibromochloromethane	BDL	5	ug/l	U
1,1,2-Trichloroethane	BDL	5	ug/l	ס
Benzene	BDL	5	ug/l	Ū
Trans-1,3-Dichloropropene	BDL	5	ug/l	ט
Bromoform	BDL	5	ug/l	Ū
4-Methyl-2-Pentanone	BDL	10	ug/1	Ū
2-Hexanone	BDL	10	ug/l	Ū
Tetrachloroethene	BDL	5	ug/1	ט
1,1,2,2-Tetrachloroethane	BDL	5	ug/l	ט
Toluene	BDL	5	ug/l	Ū
Chlorobenzene	BDL	5	ug/l	Ū

BDL

BDL

BDL



WESTON-GULF COAST LABORATORIES, INC. 2417 Bond St., University Park, Illinois 50466

Phones: (312) 534-5200 (219) 885-7077 (815) 723-753

ANALYTICAL REPORT

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater L-2-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88 Date Received: 11/18/88

GCL#:

144694

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT	
2-Chloroethyl vinyl ether	BDL	10 ug/l U	
1,3-Dichlorobenzene	BDL	10 ug/l U	_
1,2-Dichlorobenzene	BDL	10 ug/l U	
1,4-Dichlorobenzene	BDL	10 ug/l U	
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TO: Roy F. Weston, Incorporated

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

1 Weston Way

WESTON-GULF COAST LABORATORIES, INC. 2417 Bond St., University Park, Illinois 50466

Phones: (312) 534-5200 (219) 385-7077 (815) 700-7537

TAYTIL CYT BEader

NO 1989

DATE: January 6, 1989

RE: Grdwater L-3-GW-001 American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88
GCL#: 144695

VOLATILE COMPOUNDS	RESULT	DETEC	TION L	INIT
Chloromethane	BDL	10	ug/l	U
Bromomethane	BDL	10	ug/l	Ū
Vinyl Chloride	BDL	10	ug/l	O
Chloroethane	BDL	10	ug/l	Ū
Methylene Chloride	*	5	ug/1 (B
Acetone	BDL	10	ug/l	D
Carbon Disulfide	BDL .	5	ug/l	Ū
1,1-Dichloroethene	BDL	5	ug/l	Ū
1,1-Dichloroethane	BDL	5	ug/l	O
1,2-Dichloroethene (total)	BDL	5	ug/1	0
Chloroform	BDL	5	ug/l	Ū
1,2-Dichloroethane	BDL	5	ug/l	Ū
2-Butanone	BDL	10	ug/1	Ū
1,1,1-Trichloroethane	BDL	5	ug/l	O
Carbon Tetrachloride	BDL	5	ug/l	O
Vinyl Acetate	BDL	10	ug/l	U
Bromodichloromethane	BDL	5	ug/l	Ū



WESTON-GULF COAST LABORATORIES. INC. 241* Bond St., University Park, Illinois 50466

Phones: .312) 534-5290 (219) 385-707" (815) 723-753;

444_/T [4] REPSET

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater L-3-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88

VOLATILE COMPOUNDS	RESULT	DETECTION I	IMIT
1,2-Dichloropropane	BDL	5 ug/l	บ
Cis-1,3-Dichloropropene	BDL	5 ug/l	Ū
Trichloroethene	BDL	5 ug/l	U
Dibromochloromethane	BDL	5 ug/l	ט
1,1,2-Trichloroethane	BDL	5 ug/l	ט
Benzene	BDL	5 ug/l	U
Trans-1,3-Dichloropropene	BDL	5 ug/l	σ
Bromoform	BDL	5 ug/l	U
4-Methyl-2-Pentanone	BDL	10 ug/l	ט
2-Bexanone	BDL	10 ug/l	Ū
Tetrachloroethene	BDL	5 ug/l	U
1,1,2,2-Tetrachloroethane	BDL	5 ug/l	ט
Toluene	BDL	5 ug/l	U
Chlorobenzene	BDL	5 ug/l	ט
Ethylbenzene	BDL	5 ug/l	σ
Styrene	BDL	5 ug/1	Ū
Xylene	BDL	5 ug/l	U
			



WESTON-GULF COAST_ABCRATORIES, INC.
2417 Bond St., University Park, Illinois \$0466
Phones: (312) 534-5200 (219) 385-7071 (815) 123-152

7/77/2 C77 353342

TO: Roy F. Weston, Incorporated

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Grdwater L-3-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/17/88
Date Received: 11/18/88

GCL#:

144695

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U
		
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WESTON-GULF COAST LABORATORIES, INC. 2417 Bond St., University Park, Illinois 50466

Phones: (312) 534-5200 (219) 885-7077 '815) 723-7530

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water L-4-GW-001

American Home Products

W.O. # 2994-02-03

Sample Date: 11/18/88 Date Received: 11/21/88

VOLATILE COMPOUNDS	RESULT	DETEC	TION L	IMIT
Chloromethane	BDL	10	ug/1	U
Bromomethane	BDL	10	ug/l	U
Vinyl Chloride	BDL	10	ug/l	U
Chloroethane	BDL	10	ug/1	U
Methylene Chloride	75	5	ug/1	
Acetone	BDL	10	ug/l	ס
Carbon Disulfide	BDL	5	ug/1	O
1,1-Dichloroethene	BDL	5	ug/l	ס
1,1-Dichloroethane	BDL	5	ug/1	Ū
1,2-Dichloroethene (total)	BDL	5	ug/l	0
Chloroform	BDL	5	ug/1	Ū
1,2-Dichloroethane	BDL	5	ug/l	Ū
2-Butanone	BDL	10	ug/l	Ū
1,1,1-Trichloroethane	BDL	5	ug/l	Ū
Carbon Tetrachloride	BDL	5	ug/1	U
Vinyl Acetate	BDL	10	ug/l	U
Bromodichloromethane	BDL	5	ug/l	Ū



WESTON-GULF COAST LABORATORIES, INC.
2417 Bond St., University Park, Illinois 60466
Phones: (312) 534-5200 (219) 885-7077 (815) 723-752

TROPAR LAS TYLANA

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water L-4-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/18/88
Date Received: 11/21/88

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
1,2-Dichloropropane	BDL	5 ug/l U
Cis-1,3-Dichloropropene	BDL	5 ug/l U
Trichloroethene	BDL	5 ug/l U
Dibromochloromethane	BDL	5 ug/1 U
1,1,2-Trichloroethane	BDL	5 ug/1 U
Benzene	BDL	5 ug/l U
Trans-1,3-Dichloropropene	BDL	5 ug/l U
Bromoform	BDL	5 ug/l U
4-Methyl-2-Pentanone	BDL	10 ug/l U
2-Hexanone	BDL	10 ug/l 0
Tetrachloroethene	BDL	5 ug/1 U
1,1,2,2-Tetrachloroethane	BDL	5 ug/1 T
Toluene	BDL	5 ug/l U
Chlorobenzene	BDL	5 ug/l U
Ethylbenzene	BDL	5 ug/1 U
Styrene	BDL	5 ug/l U
Xylene	BDL	5 ug/l U



WESTON-GULF COAST LABORATORIES, INC. 2417 Bond St., University Park, Illinois 50466
Phones: (312) 534-5200 (219) 885-7077 (815) 723-7532

ANALYTICAL REPORT

TO: Roy F. Weston, Inc.

1 Weston Way

West Chester, PA 19380

ATTN: Ms. Sherrerd Steele

DATE: January 6, 1989

RE: Ground Water L-4-GW-001

American Home Products

W.O.# 2994-02-03

Sample Date: 11/18/88 Date Received: 11/21/88

VOLATILE COMPOUNDS	RESULT	DETECTION LIMIT
2-Chloroethyl vinyl ether	BDL	10 ug/l U
1,3-Dichlorobenzene	BDL	10 ug/l U
1,2-Dichlorobenzene	BDL	10 ug/l U
1,4-Dichlorobenzene	BDL	10 ug/l U

APPENDIX B EXPOSURE DOSE AND RISK CHARACTERIZATION TABLES



POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH INGESTION OF CHEMICALS IN RESIDENTIAL WELL WATER: UPPER UNIT EKCO HOUSEWARES MASSILLON, OHIO

					Adult			Child	
				Potential			Potential		
	Exposure Point	Cancer Slope	Chronic Reference	Chronic Daily	Potential Lifetime	Potential Hazard	Chronic Daily	Potential Lifetime	Potential Hazard
Chemical	Concentration	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	Index (HI)	Intake (CDI)	Cancer Risk	Index (HI)
	(ug/L)	(m g/kg/day) - 1	(mg/kg/day)	(mg/kg/day)	(CDIx CSF)	(CDI/(ED/70))/RfD	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD
Organics					 				
1,1-Dichloroethane	451	ND	1.0E-01	4.2E-03		1.2E~01	2.5E~03		2.9E-01
1,1-Dichloroethene	220	6.0E01	9.0E - 03	2.1E-03	1.2E - 03	6.7E~01	1.2E-03	7.2E-04	1.6E+00
1,2-Dichloroethene (total)	93.7	NC	9.0E-03	8.8E-04		2.9E ~01	5.1E-04		6.7E-01
1,1,1-Trichloroethane	5380	NC	ND	5.1E-02			2.9E ~02		
Trichloroethene	22300	1.1E-02	6.0E-03	2.1E-01	2.3E-03	1.0E+02	1.2€~01	1.3E-03	2.4E+02
Vinyl chloride	28.4	1.9E+00	ND	2.7E -04	5.1E-04		1.6E - 04	3.0E-04	
Chloroethane	3.19	NC	ND	3.0E-05			1.7E~05		
Acetone	4.25	NC	1.0E-01	4.0E-05		1.2E-03	2.3E~05		2.7E-03
1,2-Dichloroethane	9.24	9.1E-02	ND	8.7E-05	7.9E -06		5.1E~05	4.6E ~06	
2-Butanone	2.6	NC	6.0E-01	2.4E-05		1.2E-04	1.4E~05		2.8E-04
1,1,2-Trichloroethane	15.61	5.7E - 02	4.0E-03	1.5E 04	8.4E-06	1.1E-01	8.6E-05	4.9E-06	2.5E~01
Tetrachloroethene	7,71	5.2E - 02	1.0E -02	7.2E-05	3.8E-06	2.1E-02	4.2E~05	2.2E-06	4.9E-02
Toluene	12.6	NC	2.0E-01	1.2E - 04		1.7E-03	6.9E~05		4.0E-03
Ethylbenzene	1.41	NC	1.0E-01	1.3E-05		3.9E-04	7.7E-06		9.0E-04
Benzene	1.45	2.9E-02	ND	1.4E-05	3.9E −07		7.9E~06	2.3E~07	
Xylenes (total)	1.56	NC	2.0E+00	1.5E - 05		2.1E-05	8.5E~06		5.0E-05
Carbon disulfide	1.1	NC	1.0E-01	1.0E-05	-~	3.0E-04	6.0E~06		7.0E-04
Chloroform	1.99	6.1E-03	1.0E-02	1.9E-05	1.1E-07	5.5E-03	1.1E~05	6.7E-08	1.3E-02
4-Methyl-2-pentanone	2.92	NC	8.0E-02	2.7E-05		1.0E-03	1.6E-05		2.3E-03
			'	Totals:	4.1E-03	1.0E+02	Totals:	2.4E-03	2.4E+02
			!		Total Li	fetime Cancer Risk	(Adult + Child) =	6.4E-03	

CDI = (GW Concentration x IR x CF x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

Exposure Assumptions

Adult:

2 IR = Ingestion Rate (L/day)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

24 ED = Exposure Duration (years)

70 BW = Body Weight (kg)

25550 AT = Averaging Time (days)

Child:

1 IR = Ingestion Rate (L/day)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

6 ED ≈ Exposure Duration (years)

15 BW = Body Weight (kg)

TABLE B-2

POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH DERMAL ABSORPTION FROM RESIDENTIAL WELL. WATER WHILE BATHING: UPPER UNIT EKCO HOUSEWARES MASSILLON, OHIO

						Adult	j		Child	
				Chronic	Potential	Potential		Potential	Potential	
	Exposure Point	Permeability	Cancer Slope	Reference	Chronic Daily	Lifetim e	Potential Hazard	Chronic Daily	Lifetim e	Potential Hazard
Chemical	Concentration	Coefficient (PC)	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	index (Hi)	intake (CDI)	Cancer Risk	Index (HI)
	(HB/L)	(cm.hr)	(m g/kg/day) - 1	(mg/kg/day)	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD
Organics										
1,1-Dichloroethane	451	8.9E -03	ND	8.0E-02	1.1E-04		4.0E-03	4.8E-05		7.0E - 03
1,1~Dichloroethene	220	1.6E - 02	7.5E -01	7.2E -03	9.5E-05	7.1E-05	3.9E-02	4.2E-05	3.2E-05	6.8E-02
1,2-Dichloroethene (total)	93.7	1.0E-02	NC	7.2E -03	2.5E-05		1.0E-02	1.1E-05		1.8E -02
1,1,1-Trichloroethane	5380	1.7E -02	NC	ND	2.5E -03		\	1.1E-03		
Trichloroethene	22300	1.6E-02	1.4E-02	4.8E -03	9.6E - 03	1.3E-04	5.9E+00	4.3E-03	6.0E -05	1.0E+01
Vinyl chloride	28.4	7.3E-03	2.0E+00	ND	5.6E-06	1.1E-05		2.5E-06	5.0E -06	
Chloroethane	3.19	8 0E 03	NC	ND	6.9E -07			3.1E-07		
Acetone	4.25	NA	NC	8.0€ -02	0.0E+00			0.0E+00		
1,2-Dichloroethane	9.24	5.3E-03	1.1E-01	ND	1.3E-06	1.5E-07		5.9E-07	6.5E 08	
2-Butanone	2.6	1.1E-03	NC	4.8E 01	7.7E-08		4.7E-07	3.4E-08		8.3E-07
1,1,2-Trichloroethane	15.61	8.4E-03	7.1E-02	3.2E -03	3.5E-06	2.5E-07	3.2E~03	1.6E-06	1.1E-07	5.7E-03
Tetrachloroethene	7.71	4.8E-02	6.5E-02	8.0E -03	1.0E-05	6.5E-07	3.6E-03	4.4E-06	2.9E-07	6.5E-03
Toluene	12.6	4.5E-02	NÇ	1.6E -01	1.5E-05		2.8E-04	6.8E-06		5.0E - 04
Ethylbenzene	1.41	7.4E-02	NC	8.0E -02	2.8E-06		1.0E-04	1.3E-06		1.8E -04
Benzene	1.45	2.1E-02	3.6E~02	ND	8.2E-07	3.0E-08	[3.7E-07	1.3E-08	
Xylenes (total)	1.56	8.0E-02	NC	1.6E+00	3.4E-06		6.1E-06	1.5E-06		1.1E-05
Carbon disulfide	1.1	2.4E-02	NC	8.0E-02	7.1E-07		2.6E - 05	3.2E-07		4.6E-05
Chloroform	1.99	8.9E-03	7.6E -03	8.0E -03	4.8E-07	3.6E-09	1.7E 04	2.1E-07	1.6E -09	3.1E-04
4-Methyl-2-pentanone	2.92	3.3E~03	NC	6.4E-02	2.6E-07		1.2E-05	1.2E-07		2.1E-05
					Totals:	2.2E04	5.9E+00	Totals:	9.7E-05	1.1E+01
				Γ		Total Life	time Cancer Risk (Adult + Child)	32E-04	

CDI = (GW Concentration x SA x VF x CF x ET x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

NA - Not Available.

Exposure Assumptions

	Adult		Child:
23000	SA = Skin Surface Area (cm²/day)	8760	SA = Skin Surface Area (cm²/day)
0.001	VF = Volumetric Conv. Factor (L/cm³)	0.001	VF = Volumetric Conv. Factor (L/cm3)
0.001	CF = Conversion Factor (mg/ug)	0.001	CF = Conversion Factor (mg/ug)
0.25	ET = Exposure Time (hours/day)	0.25	ET = Exposure Time (hours/day)
350	EF = Exposure Frequency (days/year	350	EF = Exposure Frequency (days/year)
24	ED = Exposure Duration (years)	6	ED = Exposure Duration (years)
70	BW = Body Weight (kg)	15	BW = Body Weight (kg)
25550	AT = Averaging Time (days)	255 50	AT = Averaging Time (days)

TABLE B-3

POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH NONINGESTION (INHALATION) USES OF RESIDENTIAL WELL WATER: UPPER UNIT EKCO HOUSEWARES MASSILLON, OHIO

					Adult			Child	
,				Potential			Potential		,
' I	Exposure Point	Cancer Slope	Chronic Reference	Chronic Daily	Potential Lifetime	Potential Hazard	Chronic Daily	Potential Lifetime	Potential Hazard
Chemical	Concentration	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	Index (Hi)	Intake (CDI)	Cancer Risk	Index (HI)
	(µg/L)	(mg/kg/day) - 1	(mg/kg/day)	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/R(D	(mg/kg/day)	(CDI x CSF)	(CDV(ED/70))/RfD
Organics							 		
1,1-Dichloroethane	451	ND	1.0E-01	1.6E-02		4.6E-01	1.9E-02		2.2E+00
1,1-Dichloroethene	220	1.2E+00	9.0E-03	7.7E-03	9.3E-03	2.5E+00	9.0E - 03	1.1E-02	1.2E+01
1,2-Dichloroethene (total)	93.7	NC	9.0E-03	3.3E-03		1.1E+00	3.9E-03		5.0E+00
1,1,1-Trichloroethane	5380	NC	1.0E+00	1.9E-01		5.5E-01	2.2E - 01		2.6E+00
Trichloroethene	22300	6.0E-03	1.0E-01	7.9E-01	4.7E-03	2.3E+01	9.2E-01	5.5E-03	1.1E+02
Vinyl chloride	28.4	3.0E-01	ND:	1.0E-03	3.0E-04		1.2E-03	3.5E-04	
Chloroethane	3.19	NC	2.9E+00	1.1E-04		1.1E-04	1.3E-04		5.3E-04
Acetone	4.25	NC	1.0E-01	1.5E-04		4.4E-03	1.7E-04		2.0E - 02
1,2-Dichloroethane	9.24	9.1E-02	ND	3.3E-04	3.0E-05		3.8E-04	3.5E - 05	
2-Butanone	2.6	NC	2.9E-01	9.2E-05		9.2E-04	1.1E-04		4.3E-03
1,1,2-Trichloroethane	15.61	5.7E-02	4.0E-03	5.5E-04	3.1E-05	4.0E-01	6.4E-04	3.7E 05	1.9E+00
Tetrachioroethene	7.71	2.0E-03	1.0E-02	2.7E-04	5.4E-07	7.9E-02	3.2E-04	6.3E-07	3.7E-01
Toluene	12.6	NC	1.1E-01	4.4E-04		1.2E-02	5.2E-04		5.5E-02
Ethylbenzene	1,41	NC	2.9E-01	5.0E-05		5.0E-04	5.8E - 05		2.3E-03
Benzene	1.45	2.9E-02	ND	5.1E-05	1.5E-06		6.0E-05	1.7E - 06	
Xylenes (total)	1.56	NC	2.0E+00	5.5E-05		8.0E-05	6.4E - 05		3.7E ~04
Carbon disulfide	1.1	NC	2.9E-03	3.9E-05	~-	3.9E-02	4.5E 05	·	1.8E - 01
Chloroform	1.99	8.1E-02	1.0E-02	7.0E-05	5.7E-06	2.0E-02	8.2E-05	6.6E-06	9.5E-02
4-Methyl-2-pentanone	2.92	NC	2.0E-02	1.0E-04		1.5E-02	1.2E-04		7.0E-02
Ì				Totals:	1.4E-02	2.8E+01	Totals:	1.7E-02	1.3E+02
					Total Li	etime Cancer Risk	(Adult + Child) =	3.1E-02	

CDI = (GW Concentration x IR x K x CF x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

Exposure Assumptions

Adult:

15 IR = Inhalation Rate (m³/day)

0.5 K = Volatilization Factor (L/m³)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

24 ED = Exposure Duration (years)

70 BW = Body Weight (kg)

25550 AT = Averaging Time (days)

Child:

15 IR = Inhalation Rate (m³/day)

0.5 K = Volatilization Factor (L/m³)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

6 ED = Exposure Duration (years)

15 BW = Body Weight (kg)



POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH INGESTION OF CHEMICALS IN GARDEN PRODUCE: UPPER UMIT

EKCO HOUSEWARES

i				l.	Adult				Child		
l.	Leafy Vegetable	Garden Fruit	Root Vegetable			Potential			Potential		
1	Exposure Point	Exposure Point	Exposure Point	Cancer Stope	Chronic Reference	Chronic Daily	Potential Lifetime	Potential Hazard	Citronic Daily	Potential Lifetime	Potential Hazard
Chemical	Concentration	Concent ation	Concentration	Factor (CSF)	Dose (RfD)	intake (CDI)	Cancer Risk	Index (H)	Intake (CDI)	Cancer Risk	Index (H)
	(mg/kg)	(mg/kg)	(mg/kg)	(mg/kg/day) – 1	(mg/kg/day)	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70)) /RfD	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70)) /RfD
Organics											
1,1-Dichtroethane	4 94E-02	4.94E-02	6 90E - 01	ND	1 0E-01	1.6E- 04		4.7E-03	1.3E-04		1.5E~02
1,1-Dichbroethene	4 27E-02	4.27E-02	3.80E - 01	6.0E-01	9.0E-03	9 4E- 05	5.7E-05	3.1E-02	7.5E+ 05	4.5E-05	9.7E-02
1,2-Dichbroethene (tota)	1 32E-02	1.32E-02	1.92E-01	NC	9.0E-03	4 5E- 05		1.4E-02	3.6E- 05		4.6E-02
I,1,1- Trichbroethane	1.13E+00	1.13E+00	1.55E+01	NC	ND	3 6E-03			2.9E 03		
Ir ich broethene	1.60E+00	1.80E+00	1.31E+02	1.1E-02	6 0E-03	2.8E-02	3.1E-04	1.3E+01	2.3E- 02	2.5E-04	4.4E+01
/inyl chbride	1.21E-03	1.21E-03	2 58E-02	1.9E+00	ND	5.8E 06	1.1E- 05		4.7E-06	8.9E-06	
Chibroethane	5.96E-05	5.96E~05	3.83E-03	NC	ND	8.1E-07			6.6E-07		
Acetone	8.42E-05	8.42E-05	3 57E-03	NC	1 0E-01	7.7E-07		2.2E-05	6.3E- 07		7.3E-05
1,2- Dichtroethane	6.51E-04	8.51E-04	1.13E-02	9.1E-02	ND	2 6E- 06	2.4E-07		2.1E-06	1.9E-07	
2-Butanone	8.53E-05	8.53E-05	2 26E-03	NC	6 0E 01	5 0E- 07		2.4E-06	4.0E-07		7.9E-06
1,1,2- Trichibroethane	1.87E-03	1.87E-03	3.53E-02	5.7E-02	4.0E-03	8 OE- 06	4.5E~ 07	5.8E-03	6.4E-06	3.7E- 07	1.9E-02
letrachioroethene	2.55E-03	2.55E-03	3 01E-02	5.2E-02	1.0E-02	7.2E 06	3 7E-07	2.1E-03	5.7E-06	3.0E-07	6.7E-03
lo luene	1.97E-03	1.97E-03	4 99E-02	NC	2.0E- 01	1.1E- 05		1.6E-04	8 9E- 06		5.2E-04
Ethybenzene	1.61E-04	1.61E-04	1.12E-02	NC	1 0E-01	2.4E-08		6.9E05	1.9E 06		2.3E-04
Benzene	3.02E~04	3 02E-04	2.82E-03	2.9E-02	ND	6.9E-07	2.0E-08		5.5E- 07	1.6E-08	
(ylenes (tota)	6.30E-04	6.30E-04	1.28E~02	NC	2.0E+00	2.9E-06		4.2E-06	2.3E- 06		1.4E-05
Carbon disulfide	8.43E+04	8.43E-04	2.12E-03	NC	1.0E- Ot	7.5E-07		2.2E-05	5.6E- 07		6.5E-05
Chlorotorm	2.54E-04	2.54E-04	3.51E-03	6.1E-03	1.0E- 02	8.2E-07	5.0E-09	2.4E-04	6.6E-07	4.0E-09	7.7E-04
I-Methyl-2-pentanone	1.63E-04	1.63E-04	3.00E-03	NC	8.0E 02	6.8E~ 07		2.5E- 05	5.5E- 07		8.0E-05
					İ	Totas:	3.7E-04	1.4E+01	Totals:	3.0E-04	4.4E+01
					T		Total Lifetime	Cancer Risk (Adult +	Child) =	6.8E-04	

CDI = ((Leafy Vegetable Conc. x IR-M) + (Garden Fruit Conc. x IR-gf) + (Root Vegetable Conc. x IR-rv) x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

Exposure Assumptions

0.036 IR-W = Leafy Vegetable Ingestion Rate (lig/day)0.042 IR-gf = Garden Fruit Ingestion Rate (lig/day)

0.044 IR-rv = Root Vegetable Ingestion Rate (kg/day)

350 EF = Exposure Frequency (days/year)

24 ED = Exposure Duration (years)

70 BW = Body Weight (kg)

Adult:

25550 AT = Averaging Time (days)

Child:

0.017 IR-w = Leafy Vegetable Ingestion Rate (Ig/day)

0.028 IR-gt = Garden Fruit Ingestion Rate (lig/day)

0.031 IR-rv = Root Vegetable Ingestion Rate (kg/day)

350 EF = Exposure Frequency (days/year)

6 ED = Exposure Duration (years)

45 DW - 5-4-W-1-14

15 BW = Body Weight (kg)

TABLE B-5

POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH INGESTION OF CHEMICALS IN RESIDENTIAL WELL WATER:

LOWER UNIT EKCO HOUSEWARES

MASSILLON, OHIO

]					Adult			Child	
į				Potential			Potential		
	Exposure Point	Cancer Slope	Chronic Reference	Chronic Daily	Potential Lifetime	Potential Hazard	Chronic Daily	Potential Lifetime	Potential Hazard
Chemical	Concentration	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	Index (HI)	Intake (CDI)	Cancer Risk	Index (HI)
	(ug/L)	(m g/kg/day) - 1	(mg/kg/day)	(m g/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD	(mg/kg/day)	(CDI x CSF)	(CDV(ED/70))/RfD
Organics									·
1,1-Dichloroethane	101.67	ND	1.0E-01	9.6E-04		2.8E-02	5.6E - 04		6.5E-02
1,1-Dichloroethene	31.26	6.0E-01	9.0E - 03	2.9E-04	1.8E-04	9.5E-02	1.7E-04	1.0E-04	2.2E-01
1,2-Dichloroethene (total)	142.44	NC	9.0E - 03	1.3E-03		4.3E-01	7.8E-04		1.0E+00
1,1,1 - Trichloroethane	613.45	NC	ND	5.8E-03		1	3.4E-03		
Trichloroethene	234.19	1.1E-02	6 0E - 03	2.2E-03	2.4E-05	1.1E+00	1.3E~03	1.4E-05	2.5E+00
Vinyl chloride	8.15	1.9E+00	ND	7.7E-05	1.5E-04		4.5E-05	8.5E-05	
			L	Totals:	3.5E-04	1.6E+00	Totals:	2.0E-04	3.8E+00
į.			Γ		Total Lit	etime Cancer Risk (Adult + Child) =	5.5E-04	

CDI = (GW Concentration x IR x CF x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

Adult:

2 IR = Ingestion Rate (L/day)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

24 ED = Exposure Duration (years)

70 BW = Body Weight (kg)

25550 AT = Averaging Time (days)

Child:

Exposure Assumptions

1 IR = Ingestion Rate (L/day)

0.001 CF = Conversion Factor (mg/ug)

350 EF = Exposure Frequency (days/year)

6 ED = Exposure Duration (years)

15 BW = Body Weight (kg)



POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH DERMAL ABSORPTION FROM RESIDENTIAL WELL WATER WHILE BATHING: LOWER UNIT EKCO HOUSEWARES MASSILLON, OHIO

1						Adult			Child	
1				Chronic	Potential	Potential		Potential	Potential	
ľ	Exposure Point	Permeability	Cancer Slope	Reference	Chronic Daily	Lifetim e	Potential Hazard	Chronic Daily	Lifetime	Potential Hazard
Chemical	Concentration	Coefficient (PC)	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	Index (HI)	Intake (CDI)	Cancer Risk	Index (HI)
	(µg/L)	(cm/hr)	(mg/kg/day) - 1	(mg/kg/day)	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70))/RfD
Organics										
1,1-Dichloroethane	101.67	8.9E~03	ND	8.0E -02	2.4E-05		8.9E-04	1.1E-05		1.6E-03
1,1-Dichloroethene	31.26	1.6E -02	7.5E-01	7.2E-03	1.4E 05	1.0E-05	5.5E -03	6.0E-06	4.5E-06	9.7E-03
1,2-Dichloroethene (total)	142.44	1.0E - 02	NC	7.2E-03	3.8E-05		1.6E-02	1.7E-05		2.8E-02
1,1,1-Trichloroethane	613.45	1.7E - 02	NC	ND	2.8E-04			1.3E - 04		
Trichloroethene	234.19	1.6E-02	1.4E - 02	4.8E-03	1.0E-04	1.4E06	6.1E-02	4.5E-05	6.3E 07	1.1E-01
Vinyl chloride	8.15	7.3E-03	2.0E+00	ND	1.6E-06	3.2E-06		7.1E-07	1.4E-06	
				Į.	_Totals:	1.5E-05	8.3E-02	Totals:	6.6E-06	1.5E-01
						Total Life	ime Cancer Risk (Adult + Child):	2.1E-05	

CDI = (GW Concentration x SA x VF x CF x ET x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

Exposure Assumptions

	Adult		Child
23000	SA = Skin Surface Area (cm²/day)	8760	SA = Skin Surface Area (cm²/day)
0.001	VF = Volumetric Conv. Factor (L/cm³)	0.001	VF = Volumetric Conv. Factor (L/cm³)
0.001	CF = Conversion Factor (mg/ug)	0.001	CF = Conversion Factor (mg/ug)
0.25	ET = Exposure Time (hours/day)	0.25	ET = Exposure Time (hours/day)
350	EF = Exposure Frequency (days/year	350	EF = Exposure Frequency (days/year)
24	ED = Exposure Duration (years)	6	ED = Exposure Duration (years)
70	BW = Body Weight (kg)	15	BW = Body Weight (kg)
25550	AT = Averaging Time (days)	25550	AT = Averaging Time (days)



POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH NONINGESTION (INHALATION) USES OF RESIDENTIAL WELL WATER: LOWER UNIT EKCO HOUSEWARES MASSILLON, OHIO

1					Adult			Child_	
				Potential			Potential		
Chemical	Exposure Point Concentration (ug/L)	Cancer Slope Factor (CSF) (mg/kg/day) - 1	Chronic Reference Dose (RfD) (mg/kg/day)		Potential Lifetime Cancer Risk (CDI x CSF)	Potential Hazard Index (HI) (CDI/(ED/70))/RfD	Chronic Daily Intake (CDI) (mg/kg/day)	Potential Lifetime Cancer Risk (CDI x CSF)	Potential Hazard Index (HI) (CDI/(ED/70))/RfD
Organics									
1,1-Dichloroethane	101.67	ND	1.0E-01	3.6E-03		1.0E-01	4.2E-03		4.9E-01
1,1-Dichloroethene	31.26	1.2E+00	9.0E-03	1.1E-03	1.3E-03	3.6E-01	1.3E-03	1.5E-03	1.7E+00
1,2-Dichloroethene (total)	142.44	NC	9.0E-03	5.0E - 03		1.6E+00	5.9E~03		7.6E+00
1,1,1-Trichloroethane	613.45	NC	1.0E+00	2.2E - 02		6.3E-02	2.5E-02		2.9E~01
Trichloroethene	234.19	6.0E-03	1.0E-01	8.2E-03	4.9E-05	2 4E-01	9.6E 03	5.8E-05	1.1E+00
Vinyl chloride	8.15	3.0E-01	ND	2.9E - 04	8.6E+05		3.3E-04	1.0E-04	
				Totals:	1.5E-03	2.4E+00	Totals:	1.7E-03	1.1E+01
					Total Li	letime Cancer Risk	(Adult + Child) =	3.2E-03	

CDI = (GW Concentration x IR x K x CF x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

	Exposure Assumptions		
	Adult:		Child:
15	IR = Inhalation Rate (m³/day)	15	IR = Inhalation Rate (m³/day)
0.5	K = Volatilization Factor (L/m³)	0.5	K = Volatilization Factor (L/m³)
0.001	CF = Conversion Factor (mg/ug)	0.001	CF = Conversion Factor (mg/ug)
350	EF = Exposure Frequency (days/year)	350	EF = Exposure Frequency (days/year)
24	ED = Exposure Duration (years)	6	ED = Exposure Duration (years)
70	BW = Body Weight (kg)	15	BW = Body Weight (kg)
25550	AT = Averaging Time (days)	25550	AT = Averaging Time (days)



POTENTIAL CARCINOGENIC AND NONCARCINOGENIC HEALTH RISKS ASSOCIATED WITH INGESTION OF CHEMICALS IN GARDEN PRODUCE:

LOWER UNIT EKCO HOUSEWARES MASSILLON, OHIO

							Adult			Child	
i	Leaty Vegetable	Garden Fruit	Root Vegetable			Po tential		"	Potential		
	Exposure Point	Exposure Point	Exposure Point	Cancer Stope	Citronic Reference	Chronic Daily	Potential Lifetime	Potential Hazard	Citronic Daily	Potential Lifetime	Potential Hazard
Chemical	Concentration	Concentration	Concentration	Factor (CSF)	Dose (RfD)	Intake (CDI)	Cancer Risk	Index (H)	Intake (CDI)	Cancer Risk	Index (H)
	(ma/kg)	(mg/kg)	(mg/kg)	(mg/kg/day) – 1	(mg/kg/day)	(mg/kg/day)	(CDI x CSF)	(CDI/(ED/70)) ATD	(mg/kg/day)	(CDI x CSF)	(CDV(ED/70)) /RfD
Organics											
1,1- Dichbroethane	1.11E-02	1.11E~02	1.55E-01	ND	1.0E~ 01	3.6E-05		1.1E-03	2.9E-05		3.4E- 03
1,1- Dichbroethene	6.07E-03	6 07E-03	5 40E - 02	6 0E-01	9 OE- 03	1.3E 05	8.0E-06	4.3E-03	1.1E-05	6.4E-06	1.4E- 02
1,2- Dichbroethene (total)	2.01E-02	2 01E-02	2.92E-01	NC	9 OE - 03	6.8E 05		2.2E-02	5.4E-05		7.0E- 02
1,1,1- Trichbroethane	1.29E-01	1.29E-01	1.76E+00	NC	ND	4.1E-04			3.3E04		
Trichbro ethene	1.89E-02	1.89E-02	1.38E+00	1 1E-02	6.0E-03	2.9E-04	3 2E- 06	1.4E-01	2 4E-04	2.6E-06	4.6E-01
Vinyl chbride	3.49E~04	3 49E-04	7.40E-03	1.9E+00	ND	1.7E-06	3.1E-06		1.3E-06	2.5E-06	
1						Totab:	1.4E~05	1.7E-01	Totals:	1.2E-05	5.5E-01
Į.							Total Lifetime	Cancer Risk (Adult +	Child) =	2.6E-05	

CDI = ((Leafy Vegetable Conc, x R− N) + (Garden Fruit Conc. x R−gf) + (Root Vegetable Conc, x R+rv) x EF x ED) / (BW x AT)

NC - Not Carcinogenic.

ND - Not Determined.

	Exposure Assumptions		
	Adult:		Child:
0.036	IR- ir = Leafy Vegetable Ingestion Rate (lig/day)	0.017	IR- w = Leafy Vegetable Ingestion Rate (ig/day)
0.042	IR-gf = Garden Fruit ingestion Rate (lig/day)	0.026	IR-gf = Garden Fruit Ingestion Rate (lig/day)
0.044	IR-rv = Root Vegetable Ingestion Rate (lig/day)	0.031	IR-rv = Root Vegetable Ingestion Rate (Ing/day)
350	EF = Exposure Frequency (days/yes)	350	EF = Exposure Frequency (days/year)
24	ED = Exposure Duration (years)	6	ED = Exposure Duration (years)
70	BW ≃ Body Weight (Hg)	15	BW = Body Weight (kg)
25550	AT ≃ Averaging Time (days)	25550	AT = Averaging Time (days)

Table B-9

Central Tendency Calculations for Summary Table 6-5 in Uncertainty Analysis

Water Ingestion			Exposure *	
	Сапсег		Child	Adult
		RME ^b	5.5E-06	9.4E-06
		CT ⁶	8.5E-07	1.3E-06
		Ratio [CT/RME]	1.6E-01	1.4E - 01
	Upper	TCE RME Risk	1.3E-03	2.3E - 03
	Upper	TCE CT Risk	2.0E-04	3.1E-04
	Lower	1,1-DCE RME Risk	1.0E - 04	1.8E - 04
	Lower	1,1 - DCE CT Risk	1.6E-05	2.5E-05
	Noncancer			
		RME ^b	6.4E-05	2.7E-05
		CT ^b	3.0E-05	1.3E-05
		Ratio (CT/RME)	4.7E-01	4.7E-01
	Upper	TCE RME HQ	2.4E+02	1.0E+02
	Upper	TCE CT HQ	1.1E+02	4.7E+01
	- pp			.,
Bathing/Inhalation				
	Cancer		Child	Adult
		RME	4.1E-05	3.5E-05
		CT	9.2E-06	6.9E-06
		Ratio (CT/RME)	2.2E-01	2.0E-01
	Upper	1,1-DCE RME Risk	1.1E-02	9.3E-03
	Upper	1,1-DCE CT Risk	2.5E-03	1.8E-03
	Lower	1,1-DCE RME Risk	1.5E-03	1.3E-03
	Lower	1,1 – CT RME Risk	3.3E-04	2.5E-04
	Noncancer	•		
		RME°	4.8E - 04	1.0E - 04
		CT	3.2E - 04	6.9E - 05
		Ratio (CT/RME)	6.7E - 01	6.7E - 01
	Upper	TCE RME HQ	1.1E+02	2.3E+01
	Upper	TCE CT HQ	7.4E+01	1.5E+01
Fruit/Veg Ingestion				
	Noncancei	r		
		RME ^d	6.4E-02	1.4E - 02
		CT⁴	4.3E-02	9.2E-03
		Ratio (CT/RME)	6.7E-01	6.7E-01
	Upper	TCE RME HQ	4.4E+01	1.3E+01
	Upper	TCE CT HQ	2.9E+01	8.7E+00

^{*} These exposure values do not include the "Cw" term in the appropriate algorithms. Refer to Tables 3-1, 3-2 and 3-3 for the equations.

^b Refer to Table 3-1 and Table 6-4 to obtain variables.

 $^{^{\}circ}$ Refer to Table 3-2 and Table 6-4 to obtain variables.

^d Refer to Table 3-3 and Table 6-4 to obtain variables.

APPENDIX C GROUNDWATER DATA TREND ANALYSIS

APPENDIX C GROUNDWATER DATA TREND ANALYSIS

C.1 INTRODUCTION

Presented in this section are several discussions pertaining to the groundwater data issues which emerged following EPA Region V review of the draft baseline risk assessment.

C.2 <u>UPPER UNIT GROUNDWATER TREND ANALYSIS</u>

The following trends are apparent from an examination of the upper unit well data from sampling years 1988, 1991 and 1992:

- In the most contaminated well, D-4-30, there is a clear decrease in all VOC concentrations over the four-year period, although the 1992 concentrations of several compounds are still quite high especially 1,1-dichloroethene, 1,1,1-trichloroethane, and trichloroethane.
- In well 5-7, although VOC levels are relatively low, there is a clear decrease in VOC concentrations over the four-year period (no 1988 data was available for well 5-4, so it is not possible to draw conclusions concerning this well).
- There is a clear decrease in all VOC concentrations over the four-year period in wells I-5 and I-7.

There are no clear trends in VOC concentrations in the L-series wells, or in wells I-2 and I-4.

This brief evaluation supports the approach to use all data from 1988 to 1992 for evaluation of the upper (shallow/intermediate) unit.

C.3 WELL LOCATIONS AND YIELDS RELATIVE TO THE SHALLOW AND INTERMEDIATE GROUNDWATER UNITS

The following table identifies which EKCO wells are completed in the shallow unit and which are completed in the deep unit. It also compares the Total VOCs of the two units. The groundwater sampling data is from the most recent complete sampling round and is part of the data set used for the Risk Assessment. The data shows that VOCs in both units range from levels that are negligible to very high. While there is a wide range of VOC concentrations in both units, neither of the units demonstrates VOC concentrations that are either generally higher or lower than the other. This is demonstrated by comparing the geometric mean of the two data sets.

The following table also presents the estimated sustainable yields of the shallow and intermediate wells. These data are rough estimates, but they show that the wells are generally low producing and there is no significant difference between the units.

Shallow On-Site Wells	Total VOC March 1992 (ug/L)	Estimated Sustainable Yield (gpm)
L-1	535	1
L-2	128	1
L-3	2	<1
L-4	43	1
L-5	55	2
S-4	12	2
S-7	23	5
D-4-30	35,273	<1
I-2	1,744	5

Geometric mean-----120

Intermediate On-Site Wells	Total VOC March 1992 (ug/L)	Estimated Sustainable Yield (gpm)
I-2	1,744	5
I-4	106	5
I-5	19	5
I-7	172	2

Geometric mean -----160

C.4 R-SERIES WELLS WITH REPORT DATES 5/20/94 AND 6/22/94

The data set from 5/20/94 was a draft copy of the final data set reported on 6/22/94, and was inadvertently included in the appendix. Only data from the 6/22/94 data were evaluated in the risk assessment.

APPENDIX D RESPONSE TO EPA V COMMENTS

APPENDIX D

RESPONSE TO EPA V COMMENTS ON THE DRAFT BASELINE RISK ASSESSMENT

WESTON has reviewed the U.S. EPA Region V comments (EPA, 1994) for the draft baseline risk assessment (submitted on 11 August 1994) in the letter dated 24 October 1994 from Sally Averill, Project Manager, Technical Enforcement Section #1, RCRA Enforcement Branch. After discussing these comments with EPA Region V, we have prepared responses to each of the "General" and "Specific" numbered comments in the letter.

GENERAL COMMENTS

- 1. The comments are noted.
- a) WESTON has evaluated garden produce ingestion in the lower aquifer as recommended in the comments. Refer to Section 3 of the final report for the revisions.
 - b) <u>Dense Non-Aqueous Phase Liquids (DNAPLs</u>). DNAPLs in the upper groundwater unit underneath the EKCO facility have not been detected in two sampling studies conducted by WESTON for the RCRA Feasibility Investigation (RFI) (WESTON, 1993). The hypothetical presence of DNAPLs in unstaurated zones, composed of such chemicals as trichloroethene (TCE) and 1,1,1-trichloroethane (TCA), are unlikely based on numerous arguments which have been previously presented in this report (WESTON, 1993). However, the possibility of the presence of a DNAPL will not affect the human health risk characterization contained in this document.

WESTON conducted shallow depth-to-water (DTW) measurements on seven shallow-site wells (L-1 through L-5; S-4, and P-5) on three occasions in 1994 (15 February, 3 May and 10 August). The minimum DTW was 5.11 ft below ground surface (bgs) measured in well L-5 on 3 May. The average DTW calculated from all the shallow 1994 data was 12.50 bgs. This is based on the arithmetic mean of data collected from all seven wells during the three measurements. However, three of the wells (L-4, L-5 and S-4) included in the mean DTW estimation were in the flood plain where no houses are likely to be constructed. Exclusion of DTW data from these three wells yielded a minimum measured shallow groundwater DTW for 1994 of 14 bgs, measured in well L-2 on 3 May, and an average DTW of 17 bgs. Assuming a typical depth of 8 feet for a hypothetical residential basement, it is unlikely that measurable amounts of VOCs from hypothetical DNAPLs in the unsaturated zone would migrate to these basements during flooding, especially in view of the measured differential between the basement depth and the minimum DTW is 6 feet (i.e., minimum 14 ft bgs minus 8 ft basement depth).

- 3. "Individual" carcinogenic risks presented in the summary tables in the risk characterization chapter (Section 5) represented total individual lifetime cancer risk (adult plus child). WESTON revised the appropriate tables to reflect both child and adult risk. Appendix B presents the calculations for the child and adult exposure doses and risks.
- 4. WESTON revised Section 2 to explain the rationale for using the entire data set (1988 to 1991) for the upper unit. Also, please refer to revised Apendix C of this report, and to our response to "Specific Comment" #4 (EPA, 1994). These discussions further explain this issue, and why only 1994 data were used for the lower (bedrock) unit.

WESTON recalculated the exposure point concentrations (Section 2) and resultant doses (Section 3) and risks (Section 5) based on the inclusion of proxy concentrations that were eliminated in the draft risk assessment.

5. WESTON clarified the explanation of which on-site wells are located in the upper unit (Section 2) and more specifically, the shallow and intermediate units (Appendix C).

Appendix C in the final report also presents an evaluation of yield rates for the shallow and intermediate wells, as requested by EPA Region V, for the upper unit wells.

- 6. EKCO Housewares Facility engineering personnel confirmed that industrial (worker) use of the underlying groundwater is not used for drinking water or showering purposes. Currently, groundwater at the facility is pumped, air-stripped and then used for non-contact cooling purposes. The rationale for not evaluating workers on-site was discussed in more detail in Section 3 of the final baseline risk assessment.
- 7. WESTON recalculated inhalation doses using the method of Andelman, as described in the RAGS "Part B" guidance (EPA, 1991). The exposure algorithm (Section 3) for inhalation during showering was revised to reflect the necessary inputs for this calculation.
- 8. A discussion of central tendency issues has been included in the Uncertainty Analysis.

 Representative calculations of central tendency were presented for the key chemicals and pathways that influenced risk calculated in the "high end" (RME) approach.

SPECIFIC COMMENTS

- 1. Executive Summary, Page ES-11, Paragraph 0. DNAPLs have not been detected in two separate occasions at the EKCO Housewares Facility. Refer to the RFI and CMS reports for detailed discussions.
- 2. Section 2, Page 2-1, Paragraph 1. Because interim actions were taken on the bedrock wells up to 1994, only 1994 data were appropriate for estimating current exposure and risk. Prior to retrofitting, the overburden aquifer had significant influence on the bedrock unit. Post-IRM data are most representative of the bedrock unit. This has been clarified in Section 2. For the upper (shallow/intermediate) units, WESTON expanded the discussion on the multiple data sets used (Section 2), and evaluated groundwater trends for the VOCs in Appendix C of this report.
- 3. Table 3-4. The 2% total organic carbon (TOC) soil value was obtained from the RFI report and was based on data for the soil type characterized for the site location. This has been clarified in the text of Section 3.
- 4. Appendix A. Identical R-series well data from 5/20/94 and 6/22/94 were reevaluated as requested to determine if any errors were made. The data from 5/20 were draft forms of the data presented as "final" in the 6/22 data set. This evaluation is discussed in Appendix C.
 - Wells I-8, I-9, I-11, I-12, I-13, R-12, S-11 and S-12 are off-site well locations which were not evaluated in this risk assessment. Only on-site future uses were assessed. This is clarified in Section 2 of the final baseline risk assessment report.

REFERENCES

EPA (U.S. Environmental Protection Agency), 1989. Risk Assessment Guidance for Superfund. Vol. 1. Human Health Evaluation Manual (Part A), Interim Final. EPA/540/1-89/002.

EPA (U.S. Environmental Protection Agency), 1991. Human Health Evaluation Manual, Part B: "Development of Risk-based Preliminary Remediation Goals." OSWER Directive 9285.7-01B. 13 December 1991.

EPA (U.S. Environmental Protection Agency), 1994. Technical Review Comments on the Baseline Risk Assessment for Volatile Organic Chemicals in Groundwater at the EKCO Housewares Facility in Massillon, Ohio. Letter to P. McDonald (American Home Products Corporation) from S. Averill (U.S. EPA Region V). 24 October 1994. and Development, Washington, D.C. 10 July 1991.

WESTON (Roy F. Weston, Inc.), 1993. RCRA Facility Investigation Report, EKCO Housewares Facility, Inc., Massillon, Ohio. Roy F. Weston, Inc., 1 Weston Way, West Chester, PA 19380. May 1993